APPENDIX C

AMBIENT SAMPLING ANALYSIS (BY EVENT)

Phase I Sampling Program - 09/23/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
1,1,1-Trichloroethane	<DL	0.02	0.01	0.02	0.02	<DL	6	<DL	0.01	0.02	0.02	0.01
1,1,2,2-Tetrachloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,1,2-Trichloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,1-Dichloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,1-Dichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2,3-Trimethylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2,4-Trichlorobenzene	<DL	<DL	<DL	0.44	<DL	<DL	6	<DL	0.07	0.22	0.44	0.18
1,2,4-Trimethylbenzene	0.05	0.01	0.02	0.1	0.07	0.09	6	0.01	0.06	0.09	0.10	0.04
1,2-Dibromoethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2-Dichloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2-Dichloropropane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,3,5-Trimethylbenzene	<DL	<DL	<DL	0.05	0.02	0.02	6	<DL	0.02	0.03	0.05	0.02
1,3-Butadiene	0.02	<DL	<DL	0.06	0.07	<DL	6	<DL	0.03	0.05	0.07	0.03
1,4-Dioxane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Butanol	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Decene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Heptene	0.03	<DL	0.02	0.13	0.07	0.03	6	<DL	0.05	0.08	0.13	0.05
1-Hexene	<DL	<DL	<DL	0.03	0.02	<DL	6	<DL	0.01	0.02	0.03	0.01
1-Methylcyclohexene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Nonene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Octene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Pentene	0.07	<DL	0.04	0.16	0.17	0.07	6	<DL	0.09	0.14	0.17	0.07
1-Propanol	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Undecene	<DL	0.71	0.23	0.16	1.15	0.61	6	<DL	0.48	0.83	1.15	0.43
2,2,3-Trimethylpentane	<DL	<DL	<DL	0.04	0.08	<DL	6	<DL	0.02	0.05	0.08	0.03
2,2,4-Trimethylpentane	0.11	0.04	0.07	<DL	0.75	0.11	6	<DL	0.18	0.41	0.75	0.28
2,2,5-Trimethylhexane	<DL	<DL	<DL	0.03	0.05	<DL	6	<DL	0.01	0.03	0.05	0.02
2,3,4-Trimethylpentane	0.18	0.12	<DL	0.27	0.43	0.16	6	<DL	0.19	0.31	0.43	0.15
2,3-Dimethylbutane	0.06	0.02	0.03	0.14	0.22	0.05	6	0.02	0.09	0.15	0.22	0.08
2,3-Dimethylpentane	0.03	<DL	0.02	0.12	0.14	0.02	6	<DL	0.06	0.10	0.14	0.06
2,4,4-Trimethyl-1-Pentene	<DL	0.03	<DL	0.02	<DL	<DL	6	<DL	0.01	0.02	0.03	0.01
2,4-4-Trimethyl-2-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2,4-Dimethylpentane	0.02	<DL	0.01	0.05	0.09	0.02	6	<DL	0.03	0.06	0.09	0.03
2,5-Dimethylhexane	<DL	<DL	<DL	0.04	0.09	<DL	6	<DL	0.02	0.05	0.09	0.04
2-Ethyl-1-Butene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2-Methyl-1-Pentene	<DL	<DL	<DL	0.03	0.03	<DL	6	<DL	0.01	0.02	0.03	0.02
2-Methyl-2-Pentene	<DL	<DL	<DL	0.03	0.02	<DL	6	<DL	0.01	0.02	0.03	0.01
2-Methylheptane	0.02	<DL	0.02	0.11	0.07	0.07	6	<DL	0.05	0.08	0.11	0.04
2-Propanol	<DL	<DL	0.25	0.59	0.51	0.13	6	<DL	0.25	0.45	0.59	0.25
3-Methyl-1-Butene	<DL	<DL	<DL	<DL	0.07	<DL	6	<DL	0.01	0.04	0.07	0.03
3-Methylheptane	<DL	<DL	<DL	0.03	0.04	<DL	6	<DL	0.01	0.03	0.04	0.02
3-Methylhexane	0.12	0.1	0.13	0.27	0.27	0.18	6	0.10	0.18	0.24	0.27	0.08
3-Methylpentane	0.13	0.04	0.07	0.23	0.26	0.1	6	0.04	0.14	0.21	0.26	0.09
4-Methyl-1-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
4-Nonene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Acetaldehyde	3.37	1.74	4.84	3.9	5.2	3.92	6	1.74	3.83	4.83	5.20	1.22
Acetone (+)	4.57	3.48	6.2	4.66	5.22	4.63	6	3.48	4.79	5.52	6.20	0.89
Acetonitrile	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Acetylene	0.51	0.22	0.24	0.28	0.28	0.36	6	0.22	0.32	0.40	0.51	0.11
Acrylonitrile	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Allyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
a-Pinene	<DL	<DL	0.02	0.02	<DL	0.03	6	<DL	0.01	0.02	0.03	0.01
Benzaldehyde	0.38	0.17	0.27	0.18	0.27	0.32	6	0.17	0.27	0.33	0.38	0.08
Benzene	0.32	0.16	0.18	0.54	0.74	0.23	6	0.16	0.36	0.55	0.74	0.23
Benzyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
b-Pinene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromochloromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromodichloromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromoform	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromomethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Butyl Acrylate	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Butyraldehyde	0.22	0.22	0.49	0.68	0.33	0.49	6	0.22	0.41	0.55	0.68	0.18
c-1,2-Dichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
c-2-Butene	<DL	<DL	<DL	0.06	0.08	<DL	6	<DL	0.02	0.05	0.08	0.04
c-2-Hexene	<DL	<DL	<DL	0.01	<DL	<DL	6	<DL	NC	NC	0.01	NC
c-2-Octene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC

Phase I Sampling Program - 09/23/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
c-2-Pentene	0.05	<DL	<DL	0.12	0.13	0.03	6	<DL	0.06	0.10	0.13	0.06
c-3-Hexene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
c-3-Methyl-2-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
c-4-Methyl-2-Pentene	<DL	<DL	<DL	<DL	0.01	<DL	6	<DL	NC	NC	0.01	NC
Carbon Tetrachloride	0.11	0.13	0.11	0.11	0.11	0.11	6	0.11	0.11	0.12	0.13	0.01
Chlorobenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Chlorodifluoromethane	0.2	0.21	0.23	0.25	0.25	0.19	6	0.19	0.22	0.24	0.25	0.03
Chloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Chloroform	0.02	<DL	0.01	<DL	<DL	<DL	6	<DL	0.01	0.01	0.02	0.01
Chloromethane	0.45	0.43	0.47	0.43	0.41	0.39	6	0.39	0.43	0.45	0.47	0.03
Chloroprene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Cyclohexane	0.16	0.1	0.23	1.48	0.55	0.1	6	0.10	0.44	0.88	1.48	0.54
Cyclohexene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Cyclopentane	0.03	<DL	<DL	0.09	0.11	0.03	6	<DL	0.04	0.08	0.11	0.05
Cyclopentene	<DL	<DL	<DL	0.12	0.14	<DL	6	<DL	0.04	0.10	0.14	0.07
Dibromochloromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Dichlorodifluoromethane	0.5	0.51	0.51	0.49	0.49	0.47	6	0.47	0.50	0.51	0.51	0.02
Dichlorofluoromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Diethyl Ether	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Epichlorohydrin	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Ethane	62.3	3.3	0.66	20	11.1	8.09	6	0.66	17.58	36.29	62.30	22.93
Ethanol	1.89	1.11	1.53	2.06	1.51	1.23	6	1.11	1.56	1.85	2.06	0.37
Ethylbenzene	0.08	0.04	0.1	0.32	0.21	0.09	6	0.04	0.14	0.23	0.32	0.10
Ethylene	2.45	0.33	0.53	1.37	1.35	0.63	6	0.33	1.11	1.75	2.45	0.79
Freon 113	0.05	0.05	0.06	0.06	0.06	0.05	6	0.05	0.06	0.06	0.06	0.01
Freon 114	<DL	0.01	<DL	<DL	<DL	0.01	6	<DL	NC	NC	0.01	NC
Halocarbon 134A	0.43	0.28	0.13	0.74	0.33	0.28	6	0.13	0.37	0.53	0.74	0.21
Heptanal	<DL	0.26	0.17	0.15	0.29	0.37	6	<DL	0.21	0.31	0.37	0.13
Hexachloro-1,3-Butadiene*	<DL	<DL	<DL	0.74	<DL	<DL	6	<DL	0.12	0.37	0.74	0.30
Hexanal	0.35	0.34	0.55	0.61	0.51	0.59	6	0.34	0.49	0.59	0.61	0.12
Indan	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Indene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Isobutane	0.95	0.22	0.81	1.2	2.06	0.39	6	0.22	0.94	1.47	2.06	0.66
Isobutene + 1-Butene	0.17	0.1	0.16	0.25	0.33	0.16	6	0.10	0.20	0.26	0.33	0.08
Isobutylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Isoheptane	0.08	0.03	0.04	0.16	0.17	0.07	6	0.03	0.09	0.14	0.17	0.06
Isohexane	0.24	0.12	<DL	0.44	0.48	0.22	6	<DL	0.25	0.40	0.48	0.18
Isoprene	0.19	0.13	0.24	0.36	0.34	0.19	6	0.13	0.24	0.32	0.36	0.09
m-Dichlorobenzene	<DL	<DL	<DL	0.04	<DL	<DL	6	<DL	0.01	0.02	0.04	0.02
m-Diethylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Methyl ethyl ketone	0.44	0.37	0.67	0.54	0.6	0.59	6	0.37	0.54	0.63	0.67	0.11
Methyl t-Butylether	0.03	0.02	0.02	0.12	0.19	0.06	6	0.02	0.07	0.13	0.19	0.07
Methylcyclohexane	0.04	<DL	0.02	0.16	0.16	0.06	6	<DL	0.07	0.13	0.16	0.07
Methylcyclopentane	0.08	0.02	0.04	0.2	0.23	0.06	6	0.02	0.11	0.18	0.23	0.09
Methylene Chloride	0.05	0.03	0.13	0.04	0.08	0.03	6	0.03	0.06	0.09	0.13	0.04
Methylisobutylketone	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Naphthalene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
n-Butane	1.06	0.45	0.66	2.7	3.67	0.94	6	0.45	1.58	2.64	3.67	1.30
n-Butylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
n-Decane	0.02	<DL	0.02	0.18	0.08	0.02	6	<DL	0.05	0.11	0.18	0.07
Neohexane	0.03	<DL	<DL	0.05	0.05	0.02	6	<DL	0.03	0.04	0.05	0.02
Neopentane	<DL	<DL	<DL	0.03	0.03	<DL	6	<DL	0.01	0.02	0.03	0.02
n-Heptane	0.07	0.02	0.03	0.18	0.22	0.06	6	0.02	0.10	0.16	0.22	0.08
n-Hexane	0.13	0.06	0.11	0.29	0.34	0.13	6	0.06	0.18	0.27	0.34	0.11
n-Nonane	0.03	<DL	0.01	0.07	0.1	<DL	6	<DL	0.04	0.07	0.10	0.04
n-Octane	0.03	0.03	0.03	0.11	0.17	0.06	6	0.03	0.07	0.12	0.17	0.06
n-Pentane	0.55	0.21	0.57	1.33	1.41	0.43	6	0.21	0.75	1.16	1.41	0.50
n-Propylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
n-Undecane	0.02	<DL	<DL	0.03	0.06	0.04	6	<DL	0.03	0.04	0.06	0.02
o-Dichlorobenzene	<DL	<DL	<DL	0.07	<DL	<DL	6	<DL	0.01	0.04	0.07	0.03
p-Dichlorobenzene	0.04	0.04	<DL	0.09	0.03	0.02	6	<DL	0.04	0.06	0.09	0.03
p-Diethylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
p-Isopropyltoluene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Propane	36.6	1.58	3.59	9.15	6.06	4.4	6	1.58	10.23	20.98	36.60	13.17
Propylene	0.7	0.97	1.1	2.33	3.37	2.07	6	0.70	1.76	2.59	3.37	1.02
Styrene	0.03	<DL	0.27	0.1	0.03	<DL	6	<DL	0.07	0.16	0.27	0.10

Phase I Sampling Program - 09/23/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
t-1,2-Dichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
t-2-Butene	<DL	<DL	<DL	0.05	0.13	<DL	6	<DL	0.03	0.07	0.13	0.05
t-2-Hexene	<DL	<DL	<DL	0.02	0.02	<DL	6	<DL	0.01	0.02	0.02	0.01
t-2-Pentene	0.08	<DL	0.02	0.24	0.25	0.06	6	<DL	0.11	0.20	0.25	0.11
t-Butylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Tetrachloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Toluene	0.38	0.18	0.58	0.67	0.69	0.44	6	0.18	0.49	0.65	0.69	0.20
Trichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Trichlorofluoromethane	0.25	0.24	0.25	0.24	0.23	0.23	6	0.23	0.24	0.25	0.25	0.01
Vinyl Acetate	<DL	<DL	0.74	<DL	<DL	<DL	6	<DL	0.12	0.37	0.74	0.30
Vinyl Bromide	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Vinyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,3-Dichloropropene (Total)	ND	ND	ND	ND	ND	ND	6	<DL	NC	NC	<DL	NC
Xylene (Total)	0.36	0.15	0.31	0.94	0.68	0.47	6	0.15	0.49	0.72	0.94	0.28

<DL- Less than Detection Limit

Min- Minimum Value

Max- Maximum Value

NC- Not Calculated

STD- Standard Deviation

* Suspected Laboratory Contaminant

Phase I Sampling Program - 09/29/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
1,1,1-Trichloroethane	0.03	0.02	0.03	0.03	0.04	0.02	6	0.02	0.03	0.03	0.04	0.01
1,1,2,2-Tetrachloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,1,2-Trichloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,1-Dichloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,1-Dichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2,3-Trimethylbenzene	0.04	<DL	<DL	0.03	0.05	0.02	6	<DL	0.02	0.04	0.05	0.02
1,2,4-Trichlorobenzene	0.06	0.02	0.04	<DL	<DL	<DL	6	<DL	0.02	0.04	0.06	0.03
1,2,4-Trimethylbenzene	0.15	0.1	0.09	0.13	0.18	0.13	6	0.09	0.13	0.16	0.18	0.03
1,2-Dibromoethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2-Dichloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2-Dichloropropane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,3,5-Trimethylbenzene	0.06	0.03	0.02	0.04	0.05	0.04	6	0.02	0.04	0.05	0.06	0.01
1,3-Butadiene	0.12	0.04	0.08	0.26	0.37	0.14	6	0.04	0.17	0.27	0.37	0.12
1,4-Dioxane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Butanol	1.5	<DL	<DL	<DL	<DL	<DL	6	<DL	0.25	0.75	1.50	0.61
1-Decene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Heptene	0.41	0.12	0.16	0.2	0.27	0.12	6	0.12	0.21	0.30	0.41	0.11
1-Hexene	0.03	0.04	<DL	0.03	0.07	0.04	6	<DL	0.04	0.05	0.07	0.02
1-Methylcyclohexene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Nonene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Octene	0.1	<DL	<DL	<DL	<DL	<DL	6	<DL	0.02	0.05	0.10	0.04
1-Pentene	0.18	0.08	0.14	0.28	0.27	0.15	6	0.08	0.18	0.25	0.28	0.08
1-Propanol	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Undecene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2,2,3-Trimethylpentane	0.05	0.03	0.04	0.11	0.08	0.06	6	0.03	0.06	0.09	0.11	0.03
2,2,4-Trimethylpentane	0.37	0.21	0.4	1	0.65	0.53	6	0.21	0.53	0.75	1.00	0.28
2,2,5-Trimethylhexane	0.02	<DL	<DL	0.06	0.04	0.03	6	<DL	0.03	0.04	0.06	0.02
2,3,4-Trimethylpentane	0.27	0.19	0.34	0.62	0.38	0.34	6	0.19	0.36	0.48	0.62	0.15
2,3-Dimethylbutane	0.29	0.09	0.16	0.34	0.3	0.21	6	0.09	0.23	0.31	0.34	0.10
2,3-Dimethylpentane	0.17	0.07	0.08	0.2	0.18	0.15	6	0.07	0.14	0.19	0.20	0.05
2,4,4-Trimethyl-1-Pentene	0.04	<DL	<DL	0.08	0.02	<DL	6	<DL	0.02	0.05	0.08	0.03
2,4-4-Trimethyl-2-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2,4-Dimethylpentane	0.1	0.04	0.07	0.12	0.11	0.08	6	0.04	0.09	0.11	0.12	0.03
2,5-Dimethylhexane	0.05	0.02	0.06	0.12	0.1	0.08	6	0.02	0.07	0.10	0.12	0.04
2-Ethyl-1-Butene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2-Methyl-1-Pentene	0.03	<DL	<DL	0.04	0.04	0.02	6	<DL	0.02	0.04	0.04	0.02
2-Methyl-2-Pentene	0.06	<DL	<DL	0.05	0.06	0.02	6	<DL	0.03	0.06	0.06	0.03
2-Methylheptane	0.19	0.13	0.12	0.19	0.23	0.14	6	0.12	0.17	0.20	0.23	0.04
2-Propanol	0.56	0.34	19.8	0.65	0.51	0.48	6	0.34	3.72	10.15	19.80	7.88
3-Methyl-1-Butene	<DL	<DL	<DL	0.12	0.11	0.06	6	<DL	0.05	0.09	0.12	0.06
3-Methylheptane	0.06	0.03	<DL	0.07	0.09	0.07	6	<DL	0.05	0.08	0.09	0.03
3-Methylhexane	0.58	0.29	0.32	0.51	0.54	0.38	6	0.29	0.44	0.54	0.58	0.12
3-Methylpentane	0.61	0.23	0.24	0.6	0.59	0.39	6	0.23	0.44	0.59	0.61	0.18
4-Methyl-1-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
4-Nonene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Acetaldehyde	6.02	6.99	9.2	4.8	5.82	7.97	6	4.80	6.80	8.10	9.20	1.60
Acetone (+)	6.63	5.9	9.04	7.41	8.66	6.18	6	5.90	7.30	8.37	9.04	1.31
Acetonitrile	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Acetylene	0.91	0.47	0.49	0.59	0.61	0.64	6	0.47	0.62	0.75	0.91	0.16
Acrylonitrile	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Allyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
a-Pinene	0.12	0.09	0.09	0.06	0.07	0.05	6	0.05	0.08	0.10	0.12	0.03
Benzaldehyde	0.34	0.22	0.21	0.14	0.25	0.19	6	0.14	0.23	0.28	0.34	0.07
Benzene	0.86	0.43	0.68	1.48	0.83	0.58	6	0.43	0.81	1.11	1.48	0.36
Benzyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
b-Pinene	<DL	<DL	0.02	<DL	<DL	<DL	6	<DL	NC	NC	0.02	NC
Bromochloromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromodichloromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromoform	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromomethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Butyl Acrylate	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Butyraldehyde	0.67	0.42	0.53	0.41	0.79	0.28	6	0.28	0.52	0.67	0.79	0.19
c-1,2-Dichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
c-2-Butene	0.08	0.1	0.06	0.17	0.49	0.17	6	0.06	0.18	0.31	0.49	0.16
c-2-Hexene	<DL	<DL	<DL	<DL	0.03	0.01	6	<DL	0.01	0.02	0.03	0.01
c-2-Octene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC

Phase I Sampling Program - 09/29/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
c-2-Pentene	0.12	0.04	0.05	0.16	0.17	0.09	6	0.04	0.11	0.15	0.17	0.05
c-3-Hexene	<DL	<DL	<DL	<DL	0.05	<DL	6	<DL	0.01	0.03	0.05	0.02
c-3-Methyl-2-Pentene	0.02	<DL	<DL	<DL	0.03	<DL	6	<DL	0.01	0.02	0.03	0.01
c-4-Methyl-2-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Carbon Tetrachloride	0.14	0.13	0.12	0.16	0.12	0.12	6	0.12	0.13	0.14	0.16	0.02
Chlorobenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Chlorodifluoromethane	0.43	0.35	0.41	0.54	0.52	0.34	6	0.34	0.43	0.50	0.54	0.08
Chloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Chloroform	0.02	0.03	0.02	0.02	0.02	<DL	6	<DL	0.02	0.03	0.03	0.01
Chloromethane	0.65	0.64	0.68	0.62	0.75	0.63	6	0.62	0.66	0.70	0.75	0.05
Chloroprene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Cyclohexane	0.45	0.23	0.44	0.54	0.51	0.3	6	0.23	0.41	0.51	0.54	0.12
Cyclohexene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Cyclopentane	0.23	0.08	0.09	0.27	0.19	0.13	6	0.08	0.17	0.23	0.27	0.08
Cyclopentene	0.12	0.03	0.04	0.13	0.07	0.04	6	0.03	0.07	0.11	0.13	0.04
Dibromochloromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Dichlorodifluoromethane	0.63	0.61	0.6	0.6	0.64	0.59	6	0.59	0.61	0.63	0.64	0.02
Dichlorofluoromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Diethyl Ether	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Epichlorohydrin	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Ethane	84	17.4	17.9	36.8	19	23.5	6	17.40	33.10	54.31	84.00	25.97
Ethanol	4.98	2.34	5.41	3.83	3.73	2.82	6	2.34	3.85	4.82	5.41	1.19
Ethylbenzene	0.21	0.11	0.13	0.15	0.23	0.14	6	0.11	0.16	0.20	0.23	0.05
Ethylene	13	1.78	2.25	2.64	2.04	1.98	6	1.78	3.95	7.58	13.00	4.44
Freon 113	0.07	0.06	0.06	0.06	0.07	0.06	6	0.06	0.06	0.07	0.07	0.01
Freon 114	<DL	0.01	<DL	<DL	0.01	<DL	6	<DL	NC	NC	0.01	NC
Halocarbon 134A	0.12	0.12	0.19	0.27	0.31	0.11	6	0.11	0.19	0.26	0.31	0.09
Heptanal	0.61	0.41	0.22	0.25	0.26	0.16	6	0.16	0.32	0.45	0.61	0.17
Hexachloro-1,3-Butadiene*	0.02	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	0.02	NC
Hexanal	0.79	0.78	0.65	0.56	0.46	0.34	6	0.34	0.60	0.74	0.79	0.18
Indan	<DL	<DL	<DL	<DL	0.02	0.02	6	<DL	0.01	0.02	0.02	0.01
Indene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Isobutane	6.67	1.92	4.4	8.52	6.72	4.35	6	1.92	5.43	7.34	8.52	2.34
Isobutene + 1-Butene	0.36	0.28	0.37	0.59	1.13	0.52	6	0.28	0.54	0.79	1.13	0.31
Isobutylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Isoheptane	0.48	0.15	0.2	0.4	0.4	0.29	6	0.15	0.32	0.42	0.48	0.13
Isohexane	1.14	0.42	0.48	1.06	0.96	0.69	6	0.42	0.79	1.04	1.14	0.31
Isoprene	0.22	0.15	0.29	0.53	0.33	0.24	6	0.15	0.29	0.40	0.53	0.13
m-Dichlorobenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
m-Diethylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Methyl ethyl ketone	0.95	0.98	0.76	0.89	<DL	0.98	6	<DL	0.76	1.07	0.98	0.38
Methyl t-Butylether	0.58	1.56	0.25	0.27	0.23	0.15	6	0.15	0.51	0.94	1.56	0.54
Methylcyclohexane	0.46	0.18	0.18	0.44	0.37	0.27	6	0.18	0.32	0.42	0.46	0.13
Methylcyclopentane	0.53	0.19	0.25	0.55	0.43	0.3	6	0.19	0.38	0.50	0.55	0.15
Methylene Chloride	0.17	0.06	0.21	0.05	0.28	0.05	6	0.05	0.14	0.22	0.28	0.10
Methylisobutylketone	<DL	0.11	<DL	<DL	0.12	<DL	6	<DL	0.04	0.09	0.12	0.06
Naphthalene	<DL	<DL	<DL	<DL	0.08	0.06	6	<DL	0.02	0.05	0.08	0.04
n-Butane	11.7	2.98	4.1	11.2	8.01	6.06	6	2.98	7.34	10.30	11.70	3.62
n-Butylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
n-Decane	0.06	0.08	0.06	0.12	0.14	0.06	6	0.06	0.09	0.12	0.14	0.04
Neohexane	0.16	0.05	0.08	0.14	0.12	0.09	6	0.05	0.11	0.14	0.16	0.04
Neopentane	0.11	0.02	0.03	0.07	0.05	0.04	6	0.02	0.05	0.08	0.11	0.03
n-Heptane	0.39	0.17	0.2	0.53	0.37	0.31	6	0.17	0.33	0.44	0.53	0.13
n-Hexane	0.74	0.31	0.4	0.94	0.68	0.54	6	0.31	0.60	0.79	0.94	0.23
n-Nonane	0.07	0.07	0.08	0.14	0.18	0.1	6	0.07	0.11	0.14	0.18	0.04
n-Octane	0.17	0.12	0.15	0.3	0.25	0.17	6	0.12	0.19	0.25	0.30	0.07
n-Pentane	3.97	0.98	1.4	3.57	2.41	1.85	6	0.98	2.36	3.34	3.97	1.20
n-Propylbenzene	<DL	<DL	<DL	<DL	0.04	<DL	6	<DL	0.01	0.02	0.04	0.02
n-Undecane	0.04	0.08	0.04	0.43	0.1	0.04	6	0.04	0.12	0.25	0.43	0.15
o-Dichlorobenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
p-Dichlorobenzene	0.18	0.2	0.08	0.23	0.19	0.1	6	0.08	0.16	0.21	0.23	0.06
p-Diethylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
p-Isopropyltoluene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Propane	40.6	9.19	9.61	24.7	12.8	13.9	6	9.19	18.47	28.44	40.60	12.22
Propylene	10.4	1.6	3.08	4.07	4.78	2.7	6	1.60	4.44	6.99	10.40	3.12
Styrene	0.07	0.04	0.11	0.04	1.19	0.02	6	0.02	0.25	0.62	1.19	0.46

Phase I Sampling Program - 09/29/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
t-1,2-Dichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
t-2-Butene	0.1	0.1	0.08	0.2	0.55	0.18	6	0.08	0.20	0.35	0.55	0.18
t-2-Hexene	0.03	0.02	<DL	0.03	0.05	0.03	6	<DL	0.03	0.04	0.05	0.02
t-2-Pentene	0.24	0.06	0.11	0.32	0.32	0.18	6	0.06	0.21	0.29	0.32	0.11
t-Butylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Tetrachloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Toluene	1.37	0.63	3.67	1.01	1.42	0.83	6	0.63	1.49	2.40	3.67	1.11
Trichloroethylene	<DL	<DL	<DL	<DL	0.03	<DL	6	<DL	0.01	0.02	0.03	0.01
Trichlorofluoromethane	0.29	0.28	0.28	0.26	0.29	0.27	6	0.26	0.28	0.29	0.29	0.01
Vinyl Acetate	<DL	<DL	1.29	2.44	<DL	1.63	6	<DL	0.89	1.75	2.44	1.05
Vinyl Bromide	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Vinyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,3-Dichloropropene (Total)	ND	ND	ND	ND	ND	ND	6	<DL	NC	NC	<DL	NC
Xylene (Total)	0.9	0.5	0.54	0.64	1.05	0.6	6	0.50	0.71	0.88	1.05	0.22

<DL- Less than Detection Limit

Min- Minimum Value

Max- Maximum Value

NC- Not Calculated

STD- Standard Deviation

* Suspected Laboratory Contaminant

Phase I Sampling Program - 10/05/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
1,1,1-Trichloroethane	0.03	0.03	0.02	0.02	0.02	0.02	6	0.02	0.02	0.03	0.03	0.01
1,1,2,2-Tetrachloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,1,2-Trichloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,1-Dichloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,1-Dichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2,3-Trimethylbenzene	0.03	0.04	<DL	<DL	<DL	0.02	6	<DL	0.02	0.03	0.04	0.02
1,2,4-Trichlorobenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2,4-Trimethylbenzene	0.17	0.13	0.07	0.06	0.13	0.08	6	0.06	0.11	0.14	0.17	0.04
1,2-Dibromoethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2-Dichloroethane	0.12	0.19	0.21	0.39	0.32	0.18	6	0.12	0.24	0.32	0.39	0.10
1,2-Dichloropropane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,3,5-Trimethylbenzene	0.05	0.04	0.02	<DL	0.03	0.02	6	<DL	0.03	0.04	0.05	0.02
1,3-Butadiene	0.3	0.19	0.12	0.07	0.19	0.14	6	0.07	0.17	0.23	0.30	0.08
1,4-Dioxane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Butanol	<DL	<DL	<DL	<DL	0.32	<DL	6	<DL	0.05	0.16	0.32	0.13
1-Decene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Heptene	0.2	0.09	0.05	0.09	0.08	0.03	6	0.03	0.09	0.14	0.20	0.06
1-Hexene	0.3	0.29	0.18	0.18	0.47	0.25	6	0.18	0.28	0.37	0.47	0.11
1-Methylcyclohexene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Nonene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Octene	<DL	<DL	<DL	0.07	<DL	<DL	6	<DL	0.01	0.04	0.07	0.03
1-Pentene	0.29	0.14	0.06	0.11	0.13	0.07	6	0.06	0.13	0.20	0.29	0.08
1-Propanol	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Undecene	<DL	0.04	0.29	0.11	<DL	<DL	6	<DL	0.07	0.17	0.29	0.11
2,2,3-Trimethylpentane	0.04	0.04	<DL	0.01	0.03	0.02	6	<DL	0.02	0.04	0.04	0.02
2,2,4-Trimethylpentane	0.39	0.43	0.11	0.07	0.31	0.19	6	0.07	0.25	0.37	0.43	0.15
2,2,5-Trimethylhexane	0.02	0.02	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.02	0.01
2,3,4-Trimethylpentane	0.27	0.31	0.14	0.14	0.28	0.22	6	0.14	0.23	0.29	0.31	0.07
2,3-Dimethylbutane	0.24	0.17	0.07	0.05	0.18	0.15	6	0.05	0.14	0.20	0.24	0.07
2,3-Dimethylpentane	0.14	0.12	0.06	0.05	0.11	0.08	6	0.05	0.09	0.12	0.14	0.04
2,4,4-Trimethyl-1-Pentene	0.02	<DL	<DL	0.04	<DL	<DL	6	<DL	0.01	0.02	0.04	0.02
2,4-4-Trimethyl-2-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2,4-Dimethylpentane	0.09	0.07	0.03	0.02	0.07	0.05	6	0.02	0.06	0.08	0.09	0.03
2,5-Dimethylhexane	0.05	0.05	0.01	<DL	0.04	0.02	6	<DL	0.03	0.05	0.05	0.02
2-Ethyl-1-Butene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2-Methyl-1-Pentene	0.04	<DL	0.02	<DL	<DL	<DL	6	<DL	0.01	0.02	0.04	0.02
2-Methyl-2-Pentene	0.11	0.05	0.02	<DL	0.06	0.03	6	<DL	0.05	0.08	0.11	0.04
2-Methylheptane	0.1	0.06	0.04	0.07	0.08	0.05	6	0.04	0.07	0.08	0.10	0.02
2-Propanol	0.27	0.44	0.29	0.44	15.2	27.7	6	0.27	7.39	16.85	27.70	11.59
3-Methyl-1-Butene	0.09	<DL	<DL	<DL	<DL	<DL	6	<DL	0.02	0.05	0.09	0.04
3-Methylheptane	0.02	0.03	<DL	<DL	0.02	<DL	6	<DL	0.01	0.02	0.03	0.01
3-Methylhexane	0.4	0.31	0.21	0.19	0.36	0.23	6	0.19	0.28	0.35	0.40	0.09
3-Methylpentane	0.46	0.36	0.22	0.16	0.52	0.35	6	0.16	0.35	0.46	0.52	0.14
4-Methyl-1-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
4-Nonene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Acetaldehyde	5.61	3.13	3.78	5.09	4.95	4.38	6	3.13	4.49	5.24	5.61	0.92
Acetone (+)	4.92	4.34	4.72	3.78	5.39	6.87	6	3.78	5.00	5.87	6.87	1.06
Acetonitrile	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Acetylene	0.94	0.64	0.33	0.26	0.65	0.52	6	0.26	0.56	0.76	0.94	0.25
Acrylonitrile	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Allyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
a-Pinene	0.22	0.28	0.24	0.15	0.23	0.19	6	0.15	0.22	0.25	0.28	0.04
Benzaldehyde	0.24	<DL	0.18	0.09	<DL	0.22	6	<DL	0.12	0.21	0.24	0.11
Benzene	0.87	0.68	0.38	0.3	0.62	0.46	6	0.30	0.55	0.72	0.87	0.21
Benzyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
b-Pinene	0.03	0.07	0.05	0.02	0.01	0.03	6	0.01	0.04	0.05	0.07	0.02
Bromochloromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromodichloromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromoform	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromomethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Butyl Acrylate	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Butyraldehyde	0.45	0.16	0.2	0.16	0.33	0.24	6	0.16	0.26	0.35	0.45	0.11
c-1,2-Dichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
c-2-Butene	0.27	0.24	0.14	0.11	0.21	0.21	6	0.11	0.20	0.25	0.27	0.06
c-2-Hexene	0.03	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.03	0.01
c-2-Octene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC

Phase I Sampling Program - 10/05/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
c-2-Pentene	0.28	0.1	0.04	<DL	0.06	0.04	6	<DL	0.09	0.17	0.28	0.10
c-3-Hexene	0.04	0.02	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.04	0.02
c-3-Methyl-2-Pentene	0.04	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.04	0.02
c-4-Methyl-2-Pentene	0.05	0.01	<DL	<DL	<DL	<DL	6	<DL	0.01	0.03	0.05	0.02
Carbon Tetrachloride	0.16	0.13	0.13	0.09	0.13	0.12	6	0.09	0.13	0.15	0.16	0.02
Chlorobenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Chlorodifluoromethane	0.68	0.59	0.4	0.65	0.4	0.63	6	0.40	0.56	0.66	0.68	0.13
Chloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Chloroform	0.03	0.03	0.03	0.02	<DL	0.03	6	<DL	0.02	0.03	0.03	0.01
Chloromethane	0.7	0.72	0.62	0.53	0.72	0.68	6	0.53	0.66	0.72	0.72	0.07
Chloroprene	0.68	0.4	0.33	0.14	0.12	0.31	6	0.12	0.33	0.50	0.68	0.20
Cyclohexane	0.24	0.17	0.12	0.08	0.26	0.29	6	0.08	0.19	0.26	0.29	0.08
Cyclohexene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Cyclopentane	0.13	0.09	0.04	0.03	0.08	0.05	6	0.03	0.07	0.10	0.13	0.04
Cyclopentene	0.13	0.06	0.02	<DL	0.03	0.02	6	<DL	0.04	0.08	0.13	0.05
Dibromochloromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Dichlorodifluoromethane	0.62	0.61	0.51	0.44	0.55	0.53	6	0.44	0.54	0.60	0.62	0.07
Dichlorofluoromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Diethyl Ether	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Epichlorohydrin	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Ethane	33.9	9.97	7.76	5.23	12.1	9.23	6	5.23	13.03	21.59	33.90	10.48
Ethanol	2.84	3.15	2.36	9.34	4.6	3.74	6	2.36	4.34	6.44	9.34	2.57
Ethylbenzene	0.17	0.16	0.08	0.08	0.15	0.1	6	0.08	0.12	0.16	0.17	0.04
Ethylene	11	4.84	3.48	4.11	5.05	3.92	6	3.48	5.40	7.69	11.00	2.80
Freon 113	0.07	0.06	0.05	0.05	0.06	0.07	6	0.05	0.06	0.07	0.07	0.01
Freon 114	0.01	<DL	0.01	<DL	<DL	<DL	6	<DL	NC	NC	0.01	NC
Halocarbon 134A	0.19	0.31	0.17	0.14	0.27	0.28	6	0.14	0.23	0.28	0.31	0.07
Heptanal	0.35	<DL	<DL	<DL	<DL	<DL	6	<DL	0.06	0.18	0.35	0.14
Hexachloro-1,3-Butadiene*	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Hexanal	0.62	<DL	0.61	0.3	<DL	<DL	6	<DL	0.26	0.50	0.62	0.30
Indan	0.02	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	0.02	NC
Indene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Isobutane	3.25	2.33	1.23	0.99	2.89	1.84	6	0.99	2.09	2.82	3.25	0.90
Isobutene + 1-Butene	1.12	0.85	0.57	1.07	0.98	0.73	6	0.57	0.89	1.06	1.12	0.21
Isobutylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Isoheptane	0.28	0.25	0.11	0.1	0.25	0.14	6	0.10	0.19	0.25	0.28	0.08
Isohexane	0.86	0.54	0.31	0.23	0.71	0.49	6	0.23	0.52	0.72	0.86	0.24
Isoprene	0.26	0.38	0.19	0.1	0.23	0.2	6	0.10	0.23	0.30	0.38	0.09
m-Dichlorobenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
m-Diethylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Methyl ethyl ketone	1.07	1.02	0.96	1.34	0.97	1.29	6	0.96	1.11	1.24	1.34	0.17
Methyl t-Butylether	0.55	0.29	0.21	0.32	0.24	0.48	6	0.21	0.35	0.46	0.55	0.14
Methylcyclohexane	0.22	0.15	0.1	0.11	0.15	0.13	6	0.10	0.14	0.18	0.22	0.04
Methylcyclopentane	0.32	0.23	0.11	0.09	0.28	0.18	6	0.09	0.20	0.28	0.32	0.09
Methylene Chloride	0.07	0.06	0.05	0.71	0.08	0.18	6	0.05	0.19	0.40	0.71	0.26
Methylisobutylketone	0.17	0.15	0.2	0.39	0.1	0.18	6	0.10	0.20	0.28	0.39	0.10
Naphthalene	0.04	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.04	0.02
n-Butane	5.86	3.12	1.6	1.13	3.2	2.12	6	1.13	2.84	4.22	5.86	1.69
n-Butylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
n-Decane	0.08	0.03	0.04	0.03	0.04	0.03	6	0.03	0.04	0.06	0.08	0.02
Neohexane	0.11	0.08	0.05	0.04	0.11	0.08	6	0.04	0.08	0.10	0.11	0.03
Neopentane	0.05	0.02	0.01	<DL	0.03	0.02	6	<DL	0.02	0.04	0.05	0.02
n-Heptane	0.25	0.2	0.11	0.14	0.25	0.16	6	0.11	0.19	0.23	0.25	0.06
n-Hexane	0.46	0.61	0.39	0.46	1.15	0.79	6	0.39	0.64	0.88	1.15	0.29
n-Nonane	0.06	0.05	0.05	0.04	0.1	0.03	6	0.03	0.06	0.07	0.10	0.02
n-Octane	0.12	0.09	0.05	0.09	0.1	0.07	6	0.05	0.09	0.11	0.12	0.02
n-Pentane	2.08	1.24	0.61	0.53	1.26	0.78	6	0.53	1.08	1.56	2.08	0.58
n-Propylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
n-Undecane	0.12	0.04	0.06	0.04	0.05	0.03	6	0.03	0.06	0.08	0.12	0.03
o-Dichlorobenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
p-Dichlorobenzene	0.23	0.08	0.13	0.12	0.05	0.07	6	0.05	0.11	0.17	0.23	0.06
p-Diethylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
p-Isopropyltoluene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Propane	18	4.24	3.65	2.44	5.74	4.36	6	2.44	6.41	11.12	18.00	5.78
Propylene	3.88	2.02	1.34	1.05	1.82	1.77	6	1.05	1.98	2.79	3.88	1.00
Styrene	0.07	0.06	0.05	0.06	0.02	0.33	6	0.02	0.10	0.19	0.33	0.11

Phase I Sampling Program - 10/05/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
t-1,2-Dichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
t-2-Butene	0.31	0.27	0.17	0.1	0.24	0.26	6	0.10	0.23	0.29	0.31	0.08
t-2-Hexene	0.07	0.05	0.01	<DL	0.02	0.02	6	<DL	0.03	0.05	0.07	0.03
t-2-Pentene	0.51	0.21	0.08	<DL	0.13	0.07	6	<DL	0.17	0.32	0.51	0.18
t-Butylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Tetrachloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Toluene	1.16	1.14	0.6	0.49	1.16	0.97	6	0.49	0.92	1.17	1.16	0.30
Trichloroethylene	0.02	<DL	0.02	0.05	2.78	<DL	6	<DL	0.48	1.40	2.78	1.13
Trichlorofluoromethane	0.28	0.28	0.25	0.21	0.24	0.24	6	0.21	0.25	0.27	0.28	0.03
Vinyl Acetate	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Vinyl Bromide	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Vinyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,3-Dichloropropene (Total)	ND	ND	ND	ND	ND	ND	6	<DL	NC	NC	<DL	NC
Xylene (Total)	0.71	0.62	0.36	0.34	0.65	0.41	6	0.34	0.52	0.65	0.71	0.16

<DL- Less than Detection Limit

Min- Minimum Value

Max- Maximum Value

NC- Not Calculated

STD- Standard Deviation

* Suspected Laboratory Contaminant

Phase I Sampling Program - 10/11/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
1,1,1-Trichloroethane	0.04	INV	0.05	0.03	0.03	0.03	5	<DL	0.03	0.04	0.05	0.02
1,1,2,2-Tetrachloroethane	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
1,1,2-Trichloroethane	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
1,1-Dichloroethane	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
1,1-Dichloroethylene	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
1,2,3-Trimethylbenzene	0.02	INV	<DL	0.03	0.03	0.02	5	<DL	0.02	0.03	0.03	0.01
1,2,4-Trichlorobenzene	0.08	INV	0.08	<DL	<DL	<DL	5	<DL	0.03	0.06	0.08	0.04
1,2,4-Trimethylbenzene	0.07	INV	0.06	0.12	0.09	0.07	5	<DL	0.07	0.10	0.12	0.04
1,2-Dibromoethane	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
1,2-Dichloroethane	0.1	INV	0.13	0.14	0.13	0.14	5	<DL	0.11	0.16	0.14	0.05
1,2-Dichloropropane	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
1,3,5-Trimethylbenzene	0.03	INV	0.02	0.04	0.04	0.03	5	<DL	0.03	0.04	0.04	0.02
1,3-Butadiene	0.19	INV	0.11	0.11	0.09	0.1	5	<DL	0.10	0.15	0.19	0.06
1,4-Dioxane	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
1-Butanol	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
1-Decene	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
1-Heptene	0.03	INV	<DL	0.06	0.04	0.05	5	<DL	0.03	0.05	0.06	0.03
1-Hexene	0.03	INV	0.03	0.06	0.04	0.04	5	<DL	0.03	0.05	0.06	0.02
1-Methylcyclohexene	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
1-Nonene	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
1-Octene	0.03	INV	<DL	<DL	<DL	<DL	5	<DL	0.01	0.02	0.03	0.01
1-Pentene	0.05	INV	0.05	0.11	0.06	0.08	5	<DL	0.06	0.09	0.11	0.04
1-Propanol	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
1-Undecene	0.23	INV	0.06	0.09	0.12	<DL	5	<DL	0.08	0.16	0.23	0.09
2,2,3-Trimethylpentane	0.02	INV	0.01	0.03	0.02	0.02	5	<DL	0.02	0.03	0.03	0.01
2,2,4-Trimethylpentane	0.17	INV	0.15	0.28	0.21	0.18	5	<DL	0.17	0.25	0.28	0.09
2,2,5-Trimethylhexane	0.02	INV	0.02	0.02	0.02	0.02	5	<DL	0.02	0.02	0.02	0.01
2,3,4-Trimethylpentane	0.12	INV	0.12	0.18	0.15	0.14	5	<DL	0.12	0.17	0.18	0.06
2,3-Dimethylbutane	0.11	INV	0.08	0.12	0.08	0.08	5	<DL	0.08	0.12	0.12	0.04
2,3-Dimethylpentane	0.07	INV	0.06	0.09	0.07	0.04	5	<DL	0.06	0.08	0.09	0.03
2,4,4-Trimethyl-1-Pentene	0.01	INV	<DL	0.02	0.03	<DL	5	<DL	0.01	0.02	0.03	0.01
2,4-4-Trimethyl-2-Pentene	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
2,4-Dimethylpentane	0.04	INV	0.03	0.05	0.03	0.03	5	<DL	0.03	0.04	0.05	0.02
2,5-Dimethylhexane	0.02	INV	0.02	0.03	0.02	0.01	5	<DL	0.02	0.03	0.03	0.01
2-Ethyl-1-Butene	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
2-Methyl-1-Pentene	<DL	INV	<DL	0.02	<DL	<DL	5	<DL	NC	NC	0.02	NC
2-Methyl-2-Pentene	<DL	INV	<DL	0.02	0.01	<DL	5	<DL	0.01	0.01	0.02	0.01
2-Methylheptane	0.04	INV	0.02	0.07	0.05	0.05	5	<DL	0.04	0.06	0.07	0.02
2-Propanol	<DL	INV	0.69	0.29	0.5	0.52	5	<DL	0.33	0.59	0.69	0.29
3-Methyl-1-Butene	0.05	INV	<DL	<DL	<DL	<DL	5	<DL	0.01	0.03	0.05	0.02
3-Methylheptane	0.02	INV	<DL	0.02	<DL	<DL	5	<DL	0.01	0.02	0.02	0.01
3-Methylhexane	0.17	INV	0.16	0.26	0.18	0.19	5	<DL	0.16	0.24	0.26	0.09
3-Methylpentane	0.22	INV	0.19	0.24	0.22	0.18	5	<DL	0.18	0.25	0.24	0.09
4-Methyl-1-Pentene	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
4-Nonene	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
Acetaldehyde	2.51	INV	2.44	4.84	2.48	7.77	5	<DL	3.34	5.72	7.77	2.66
Acetone (+)	3.59	INV	6.08	7.2	4.75	5.74	5	<DL	4.56	6.84	7.20	2.55
Acetonitrile	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
Acetylene	0.6	INV	0.55	0.7	0.64	0.63	5	<DL	0.52	0.75	0.70	0.26
Acrylonitrile	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
Allyl Chloride	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
a-Pinene	0.07	INV	0.08	0.08	0.17	0.06	5	<DL	0.08	0.13	0.17	0.05
Benzaldehyde	0.27	INV	0.33	0.2	0.11	0.61	5	<DL	0.25	0.44	0.61	0.21
Benzene	0.5	INV	0.51	0.77	0.6	0.81	5	<DL	0.53	0.79	0.81	0.29
Benzyl Chloride	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
b-Pinene	<DL	INV	0.02	0.02	<DL	<DL	5	<DL	0.01	0.02	0.02	0.01
Bromochloromethane	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
Bromodichloromethane	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
Bromoform	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
Bromomethane	0.03	INV	0.03	0.02	<DL	0.02	5	<DL	0.02	0.03	0.03	0.01
Butyl Acrylate	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
Butyraldehyde	0.19	INV	0.13	0.42	0.07	0.31	5	<DL	0.19	0.33	0.42	0.16
c-1,2-Dichloroethylene	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
c-2-Butene	0.07	INV	0.02	0.17	0.19	1.77	5	<DL	0.37	0.99	1.77	0.69
c-2-Hexene	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
c-2-Octene	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC

Phase I Sampling Program - 10/11/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
c-2-Pentene	0.02	INV	<DL	0.04	0.02	0.02	5	<DL	0.02	0.03	0.04	0.02
c-3-Hexene	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
c-3-Methyl-2-Pentene	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
c-4-Methyl-2-Pentene	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
Carbon Tetrachloride	0.16	INV	0.17	0.18	0.14	0.16	5	<DL	0.14	0.20	0.18	0.07
Chlorobenzene	0.02	INV	0.01	0.02	0.01	0.02	5	<DL	0.01	0.02	0.02	0.01
Chlorodifluoromethane	0.47	INV	0.41	0.43	0.35	0.88	5	<DL	0.42	0.67	0.88	0.28
Chloroethane	0.08	INV	<DL	<DL	<DL	<DL	5	<DL	0.01	0.04	0.08	0.03
Chloroform	0.03	INV	0.04	0.04	0.03	0.04	5	<DL	0.03	0.04	0.04	0.02
Chloromethane	0.5	INV	0.57	0.55	0.47	0.51	5	<DL	0.43	0.63	0.57	0.22
Chloroprene	0.3	INV	0.34	0.28	0.19	0.26	5	<DL	0.23	0.34	0.34	0.12
Cyclohexane	0.11	INV	0.1	0.11	0.15	0.13	5	<DL	0.10	0.15	0.15	0.05
Cyclohexene	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
Cyclopentane	0.05	INV	0.04	0.06	0.08	0.06	5	<DL	0.05	0.07	0.08	0.03
Cyclopentene	<DL	INV	<DL	0.02	<DL	0.02	5	<DL	0.01	0.02	0.02	0.01
Dibromochloromethane	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
Dichlorodifluoromethane	0.6	INV	0.62	0.59	0.5	0.54	5	<DL	0.48	0.69	0.62	0.24
Dichlorofluoromethane	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
Diethyl Ether	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
Epichlorohydrin	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
Ethane	5.95	INV	6.19	6.75	6.74	6.72	5	<DL	5.39	7.77	6.75	2.66
Ethanol	1.61	INV	1.34	1.31	1.58	1.19	5	<DL	1.17	1.71	1.61	0.60
Ethylbenzene	0.11	INV	0.13	0.25	0.21	0.26	5	<DL	0.16	0.25	0.26	0.10
Ethylene	1.51	INV	1.65	2.21	1.89	1.99	5	<DL	1.54	2.25	2.21	0.79
Freon 113	0.1	INV	0.08	0.07	0.06	0.06	5	<DL	0.06	0.09	0.10	0.03
Freon 114	0.02	INV	0.02	0.01	0.01	0.01	5	<DL	0.01	0.02	0.02	0.01
Halocarbon 134A	0.1	INV	0.21	0.09	0.12	0.08	5	<DL	0.10	0.16	0.21	0.07
Heptanal	<DL	INV	<DL	0.19	<DL	<DL	5	<DL	0.03	0.10	0.19	0.08
Hexachloro-1,3-Butadiene*	0.04	INV	0.05	<DL	<DL	<DL	5	<DL	0.02	0.04	0.05	0.02
Hexanal	0.22	INV	0.19	0.5	0.24	0.4	5	<DL	0.26	0.41	0.50	0.17
Indan	<DL	INV	<DL	0.02	<DL	0.02	5	<DL	0.01	0.02	0.02	0.01
Indene	<DL	INV	<DL	<DL	<DL	0.01	5	<DL	NC	NC	0.01	NC
Isobutane	1.32	INV	1.43	1.47	1.52	1.37	5	<DL	1.19	1.71	1.52	0.58
Isobutene + 1-Butene	0.16	INV	0.15	0.47	0.37	1.19	5	<DL	0.39	0.77	1.19	0.43
Isobutylbenzene	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
Isoheptane	0.11	INV	0.1	0.11	0.13	0.1	5	<DL	0.09	0.13	0.13	0.05
Isohexane	0.31	INV	0.29	0.41	0.31	0.33	5	<DL	0.28	0.40	0.41	0.14
Isoprene	0.21	INV	0.23	0.26	0.15	0.09	5	<DL	0.16	0.24	0.26	0.10
m-Dichlorobenzene	0.02	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	0.02	NC
m-Diethylbenzene	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
Methyl ethyl ketone	0.52	INV	0.63	1.66	1.64	6.71	5	<DL	1.86	4.06	6.71	2.46
Methyl t-Butylether	0.06	INV	0.05	0.45	0.22	0.61	5	<DL	0.23	0.45	0.61	0.25
Methylcyclohexane	0.07	INV	0.1	0.1	0.1	0.1	5	<DL	0.08	0.11	0.10	0.04
Methylcyclopentane	0.11	INV	0.1	0.15	0.13	0.13	5	<DL	0.10	0.15	0.15	0.05
Methylene Chloride	0.14	INV	0.13	0.19	0.36	1.88	5	<DL	0.45	1.09	1.88	0.71
Methylisobutylketone	<DL	INV	<DL	0.1	<DL	<DL	5	<DL	0.02	0.05	0.10	0.04
Naphthalene	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
n-Butane	2.6	INV	2.44	2.92	2.64	4.97	5	<DL	2.60	4.01	4.97	1.58
n-Butylbenzene	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
n-Decane	0.03	INV	0.03	0.03	0.08	0.03	5	<DL	0.03	0.06	0.08	0.03
Neohexane	0.07	INV	0.07	0.09	0.09	0.07	5	<DL	0.07	0.09	0.09	0.03
Neopentane	0.02	INV	0.02	0.03	0.02	0.02	5	<DL	0.02	0.03	0.03	0.01
n-Heptane	0.1	INV	0.1	0.13	0.13	0.09	5	<DL	0.09	0.13	0.13	0.05
n-Hexane	0.31	INV	0.33	0.39	0.36	0.42	5	<DL	0.30	0.44	0.42	0.15
n-Nonane	0.04	INV	0.03	0.04	0.09	0.03	5	<DL	0.04	0.06	0.09	0.03
n-Octane	0.05	INV	0.06	0.09	0.09	0.08	5	<DL	0.06	0.09	0.09	0.03
n-Pentane	0.87	INV	0.93	0.97	2.58	1.12	5	<DL	1.08	1.83	2.58	0.84
n-Propylbenzene	<DL	INV	<DL	0.04	<DL	<DL	5	<DL	0.01	0.02	0.04	0.02
n-Undecane	0.03	INV	0.04	0.03	0.07	0.03	5	<DL	0.03	0.05	0.07	0.02
o-Dichlorobenzene	0.02	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	0.02	NC
p-Dichlorobenzene	0.13	INV	0.58	0.13	0.05	0.02	5	<DL	0.15	0.35	0.58	0.22
p-Diethylbenzene	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
p-Isopropyltoluene	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
Propane	3.16	INV	4	4.02	3.5	3.86	5	<DL	3.09	4.48	4.02	1.55
Propylene	0.84	INV	1.05	1.4	1.4	1.45	5	<DL	1.02	1.52	1.45	0.56
Styrene	0.06	INV	0.26	0.12	0.95	0.28	5	<DL	0.28	0.59	0.95	0.35

Phase I Sampling Program - 10/11/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
t-1,2-Dichloroethylene	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
t-2-Butene	0.07	INV	<DL	0.2	0.24	2.33	5	<DL	0.47	1.29	2.33	0.92
t-2-Hexene	<DL	INV	<DL	0.02	0.01	<DL	5	<DL	0.01	0.01	0.02	0.01
t-2-Pentene	0.03	INV	0.02	0.07	0.05	0.03	5	<DL	0.03	0.05	0.07	0.02
t-Butylbenzene	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
Tetrachloroethylene	0.01	INV	0.02	<DL	<DL	<DL	5	<DL	0.01	0.01	0.02	0.01
Toluene	0.62	INV	0.69	1.08	1.38	0.94	5	<DL	0.79	1.21	1.38	0.47
Trichloroethylene	0.04	INV	0.12	0.02	0.07	<DL	5	<DL	0.04	0.08	0.12	0.05
Trichlorofluoromethane	0.4	INV	0.33	0.31	0.26	0.29	5	<DL	0.27	0.39	0.40	0.14
Vinyl Acetate	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
Vinyl Bromide	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
Vinyl Chloride	<DL	INV	<DL	<DL	<DL	<DL	5	<DL	NC	NC	<DL	NC
1,3-Dichloropropene (Total)	ND	INV	ND	ND	ND	ND	5	<DL	NC	NC	<DL	NC
Xylene (Total)	0.31	INV	0.3	0.65	0.49	0.5	5	<DL	0.38	0.58	0.65	0.23

<DL- Less than Detection Limit

INV- Data Invalidated

Min- Minimum Value

Max- Maximum Value

NC- Not Calculated

STD- Standard Deviation

* Suspected Laboratory Contaminant

Phase I Sampling Program - 10/17/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
1,1,1-Trichloroethane	0.03	0.03	0.03	<DL	0.03	0.02	6	<DL	0.02	0.03	0.03	0.01
1,1,2,2-Tetrachloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,1,2-Trichloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,1-Dichloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,1-Dichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2,3-Trimethylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2,4-Trichlorobenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2,4-Trimethylbenzene	0.11	0.04	0.06	0.08	0.1	0.05	6	0.04	0.07	0.10	0.11	0.03
1,2-Dibromoethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2-Dichloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2-Dichloropropane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,3,5-Trimethylbenzene	0.03	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.03	0.01
1,3-Butadiene	0.13	0.05	0.09	0.22	0.27	0.18	6	0.05	0.16	0.22	0.27	0.08
1,4-Dioxane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Butanol	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Decene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Heptene	0.18	<DL	<DL	<DL	<DL	<DL	6	<DL	0.03	0.09	0.18	0.07
1-Hexene	<DL	<DL	0.05	<DL	0.09	<DL	6	<DL	0.02	0.05	0.09	0.04
1-Methylcyclohexene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Nonene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Octene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Pentene	0.24	0.16	0.27	0.29	0.39	0.27	6	0.16	0.27	0.33	0.39	0.07
1-Propanol	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Undecene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2,2,3-Trimethylpentane	0.03	0.02	0.02	0.03	0.03	<DL	6	<DL	0.02	0.03	0.03	0.01
2,2,4-Trimethylpentane	0.43	0.24	0.31	0.44	0.34	0.24	6	0.24	0.33	0.41	0.44	0.09
2,2,5-Trimethylhexane	0.02	<DL	<DL	<DL	0.02	<DL	6	<DL	0.01	0.02	0.02	0.01
2,3,4-Trimethylpentane	0.26	0.2	0.26	0.38	0.23	0.29	6	0.20	0.27	0.32	0.38	0.06
2,3-Dimethylbutane	0.27	0.11	0.16	0.21	0.23	0.13	6	0.11	0.19	0.24	0.27	0.06
2,3-Dimethylpentane	0.16	0.06	0.08	0.11	0.13	0.07	6	0.06	0.10	0.13	0.16	0.04
2,4,4-Trimethyl-1-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2,4-4-Trimethyl-2-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2,4-Dimethylpentane	0.11	<DL	<DL	0.1	0.09	<DL	6	<DL	0.05	0.10	0.11	0.06
2,5-Dimethylhexane	0.03	0.01	0.02	0.03	0.03	<DL	6	<DL	0.02	0.03	0.03	0.01
2-Ethyl-1-Butene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2-Methyl-1-Pentene	0.04	<DL	<DL	<DL	0.07	<DL	6	<DL	0.02	0.04	0.07	0.03
2-Methyl-2-Pentene	0.04	<DL	0.04	<DL	0.07	<DL	6	<DL	0.03	0.05	0.07	0.03
2-Methylheptane	0.1	0.04	0.06	0.07	0.09	0.07	6	0.04	0.07	0.09	0.10	0.02
2-Propanol	<DL	4.7	2.04	6.13	<DL	1.98	6	<DL	2.48	4.51	6.13	2.49
3-Methyl-1-Butene	0.12	<DL	0.11	0.14	0.19	0.1	6	<DL	0.11	0.16	0.19	0.06
3-Methylheptane	0.04	<DL	<DL	<DL	0.03	<DL	6	<DL	0.01	0.03	0.04	0.02
3-Methylhexane	0.3	0.09	0.15	0.19	0.25	0.17	6	0.09	0.19	0.25	0.30	0.07
3-Methylpentane	0.46	0.13	0.2	0.26	0.44	0.2	6	0.13	0.28	0.39	0.46	0.14
4-Methyl-1-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
4-Nonene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Acetaldehyde	1.35	2.71	2.59	2.54	2.6	3.35	6	1.35	2.52	3.05	3.35	0.65
Acetone (+)	1.87	2.25	0.62	7.54	1.03	2.62	6	0.62	2.66	4.70	7.54	2.51
Acetonitrile	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Acetylene	1.14	0.46	0.75	0.85	0.87	0.58	6	0.46	0.78	0.97	1.14	0.24
Acrylonitrile	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Allyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
a-Pinene	<DL	0.02	0.03	0.02	<DL	<DL	6	<DL	0.01	0.02	0.03	0.01
Benzaldehyde	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Benzene	0.86	0.26	0.42	0.78	0.65	0.62	6	0.26	0.60	0.78	0.86	0.22
Benzyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
b-Pinene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromochloromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromodichloromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromoform	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromomethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Butyl Acrylate	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Butyraldehyde	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
c-1,2-Dichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
c-2-Butene	0.16	0.16	0.3	0.29	1.6	0.27	6	0.16	0.46	0.92	1.60	0.56
c-2-Hexene	<DL	<DL	<DL	<DL	0.03	<DL	6	<DL	0.01	0.02	0.03	0.01
c-2-Octene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC

Phase I Sampling Program - 10/17/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
c-2-Pentene	0.14	0.08	0.12	0.11	0.2	0.05	6	0.05	0.12	0.16	0.20	0.05
c-3-Hexene	0.04	<DL	0.01	<DL	0.06	<DL	6	<DL	0.02	0.04	0.06	0.03
c-3-Methyl-2-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
c-4-Methyl-2-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Carbon Tetrachloride	0.11	0.12	0.12	0.11	0.11	0.11	6	0.11	0.11	0.12	0.12	0.01
Chlorobenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Chlorodifluoromethane	0.46	1.2	1.7	2.4	0.47	1.14	6	0.46	1.23	1.84	2.40	0.74
Chloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Chloroform	0.03	0.02	0.02	<DL	0.06	<DL	6	<DL	0.02	0.04	0.06	0.02
Chloromethane	0.53	0.58	0.57	0.55	0.46	0.45	6	0.45	0.52	0.57	0.58	0.06
Chloroprene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Cyclohexane	0.22	0.04	0.07	0.14	0.1	0.08	6	0.04	0.11	0.16	0.22	0.06
Cyclohexene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Cyclopentane	0.14	<DL	0.06	0.09	0.1	0.06	6	<DL	0.08	0.11	0.14	0.05
Cyclopentene	0.08	<DL	0.04	<DL	0.05	<DL	6	<DL	0.03	0.06	0.08	0.03
Dibromochloromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Dichlorodifluoromethane	0.6	0.64	0.75	0.65	0.66	0.59	6	0.59	0.65	0.69	0.75	0.06
Dichlorofluoromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Diethyl Ether	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Epichlorohydrin	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Ethane	19.2	4.34	7.78	11.8	8.19	10.1	6	4.34	10.24	14.36	19.20	5.06
Ethanol	<DL	10.9	3.56	2.09	2.7	4.71	6	<DL	3.99	7.04	10.90	3.73
Ethylbenzene	0.13	0.05	0.12	0.1	0.13	0.09	6	0.05	0.10	0.13	0.13	0.03
Ethylene	6.1	1.18	2.07	3.76	2.64	3.16	6	1.18	3.15	4.54	6.10	1.70
Freon 113	0.07	0.07	0.07	0.07	0.07	0.06	6	0.06	0.07	0.07	0.07	0.00
Freon 114	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Halocarbon 134A	0.15	0.78	0.49	0.28	0.1	0.43	6	0.10	0.37	0.58	0.78	0.25
Heptanal	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Hexachloro-1,3-Butadiene*	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Hexanal	<DL	0.16	<DL	<DL	<DL	<DL	6	<DL	0.03	0.08	0.16	0.07
Indan	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Indene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Isobutane	4.16	1.5	2.12	6.98	6.54	2.46	6	1.50	3.96	5.87	6.98	2.35
Isobutene + 1-Butene	0.33	0.18	0.28	0.51	2.77	0.64	6	0.18	0.79	1.59	2.77	0.99
Isobutylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Isoheptane	0.3	0.1	0.15	0.2	0.25	0.16	6	0.10	0.19	0.25	0.30	0.07
Isohexane	1.12	0.29	0.53	0.64	1.22	0.43	6	0.29	0.71	1.01	1.22	0.38
Isoprene	0.09	0.09	0.09	0.1	0.07	0.09	6	0.07	0.09	0.10	0.10	0.01
m-Dichlorobenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
m-Diethylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Methyl ethyl ketone	<DL	<DL	<DL	<DL	0.64	<DL	6	<DL	0.11	0.32	0.64	0.26
Methyl t-Butylether	0.24	<DL	0.03	<DL	1.79	0.29	6	<DL	0.39	0.96	1.79	0.70
Methylcyclohexane	0.32	0.1	0.17	0.16	0.11	0.13	6	0.10	0.17	0.23	0.32	0.08
Methylcyclopentane	0.36	0.09	0.15	0.2	0.25	0.13	6	0.09	0.20	0.28	0.36	0.10
Methylene Chloride	0.21	1.45	1.81	0.45	0.11	1.17	6	0.11	0.87	1.44	1.81	0.71
Methylisobutylketone	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Naphthalene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
n-Butane	12.1	4.67	7.06	6.43	18.5	4.27	6	4.27	8.84	13.33	18.50	5.50
n-Butylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
n-Decane	0.02	0.02	0.03	0.04	0.03	<DL	6	<DL	0.02	0.03	0.04	0.01
Neohexane	0.09	0.03	0.04	0.05	0.03	<DL	6	<DL	0.04	0.06	0.09	0.03
Neopentane	0.08	0.02	0.02	0.02	0.05	0.01	6	0.01	0.03	0.06	0.08	0.03
n-Heptane	0.26	0.06	0.1	0.11	0.12	0.08	6	0.06	0.12	0.18	0.26	0.07
n-Hexane	0.53	0.14	0.25	0.26	0.43	0.2	6	0.14	0.30	0.42	0.53	0.15
n-Nonane	0.07	0.04	0.05	0.08	0.07	0.06	6	0.04	0.06	0.07	0.08	0.01
n-Octane	0.1	0.05	0.07	0.08	0.09	0.08	6	0.05	0.08	0.09	0.10	0.02
n-Pentane	3.78	0.69	1.28	1.64	2.14	1.3	6	0.69	1.81	2.69	3.78	1.08
n-Propylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
n-Undecane	0.02	0.01	0.02	<DL	0.02	0.02	6	<DL	0.02	0.02	0.02	0.01
o-Dichlorobenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
p-Dichlorobenzene	0.09	0.82	0.35	0.16	0.08	0.58	6	0.08	0.35	0.59	0.82	0.30
p-Diethylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
p-Isopropyltoluene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Propane	13.3	5.76	9.55	14.7	5.84	8.65	6	5.76	9.63	12.68	14.70	3.73
Propylene	5.93	3.04	3.21	12.3	3.07	8.48	6	3.04	6.01	9.08	12.30	3.77
Styrene	0.04	0.03	0.09	0.32	<DL	<DL	6	<DL	0.08	0.18	0.32	0.12

Phase I Sampling Program - 10/17/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
t-1,2-Dichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
t-2-Butene	0.16	0.2	0.29	0.38	1.9	0.47	6	0.16	0.57	1.11	1.90	0.66
t-2-Hexene	0.04	0.01	0.02	<DL	0.08	<DL	6	<DL	0.03	0.05	0.08	0.03
t-2-Pentene	0.35	0.13	0.25	0.21	0.43	0.12	6	0.12	0.25	0.35	0.43	0.12
t-Butylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Tetrachloroethylene	0.02	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	0.02	NC
Toluene	1.43	0.56	0.96	1.38	1.21	0.89	6	0.56	1.07	1.34	1.43	0.33
Trichloroethylene	0.09	0.82	0.4	0.17	0.03	0.44	6	0.03	0.33	0.56	0.82	0.29
Trichlorofluoromethane	0.29	0.35	0.34	0.33	0.29	0.33	6	0.29	0.32	0.34	0.35	0.03
Vinyl Acetate	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Vinyl Bromide	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Vinyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,3-Dichloropropene (Total)	ND	ND	ND	ND	ND	ND	6	<DL	NC	NC	<DL	NC
Xylene (Total)	0.59	0.19	0.51	0.45	0.58	0.35	6	0.19	0.45	0.57	0.59	0.15

<DL- Less than Detection Limit

Min- Minimum Value

Max- Maximum Value

NC- Not Calculated

STD- Standard Deviation

* Suspected Laboratory Contaminant

Phase I Sampling Program - 10/23/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
1,1,1-Trichloroethane	0.04	0.03	0.02	0.03	0.04	0.05	6	0.02	0.04	0.04	0.05	0.01
1,1,2,2-Tetrachloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,1,2-Trichloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,1-Dichloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,1-Dichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2,3-Trimethylbenzene	0.03	<DL	<DL	0.02	0.02	0.01	6	<DL	0.01	0.02	0.03	0.01
1,2,4-Trichlorobenzene	0.09	0.02	0.05	<DL	<DL	<DL	6	<DL	0.03	0.06	0.09	0.04
1,2,4-Trimethylbenzene	0.12	0.02	0.07	0.1	0.07	0.06	6	0.02	0.07	0.10	0.12	0.03
1,2-Dibromoethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2-Dichloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2-Dichloropropane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,3,5-Trimethylbenzene	0.04	<DL	0.02	0.03	0.02	0.02	6	<DL	0.02	0.03	0.04	0.01
1,3-Butadiene	0.13	0.02	0.05	0.24	0.16	0.09	6	0.02	0.12	0.18	0.24	0.08
1,4-Dioxane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Butanol	<DL	<DL	0.25	<DL	<DL	0.23	6	<DL	0.08	0.18	0.25	0.12
1-Decene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Heptene	0.11	0.07	0.14	0.27	0.24	0.2	6	0.07	0.17	0.24	0.27	0.08
1-Hexene	0.07	<DL	0.04	0.03	0.04	0.02	6	<DL	0.03	0.05	0.07	0.02
1-Methylcyclohexene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Nonene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Octene	<DL	0.04	0.06	<DL	<DL	<DL	6	<DL	0.02	0.04	0.06	0.03
1-Pentene	0.16	0.09	0.1	0.2	0.2	0.12	6	0.09	0.15	0.18	0.20	0.05
1-Propanol	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Undecene	0.22	0.26	0.37	<DL	<DL	<DL	6	<DL	0.14	0.27	0.37	0.16
2,2,3-Trimethylpentane	0.03	0.01	0.02	0.08	0.04	0.03	6	0.01	0.04	0.05	0.08	0.02
2,2,4-Trimethylpentane	0.23	0.06	0.12	0.7	0.3	0.16	6	0.06	0.26	0.45	0.70	0.23
2,2,5-Trimethylhexane	0.02	<DL	0.01	0.03	0.02	0.01	6	<DL	0.02	0.02	0.03	0.01
2,3,4-Trimethylpentane	0.16	<DL	<DL	0.35	0.21	<DL	6	<DL	0.12	0.24	0.35	0.15
2,3-Dimethylbutane	0.19	0.06	0.11	0.34	0.21	0.13	6	0.06	0.17	0.25	0.34	0.10
2,3-Dimethylpentane	0.07	0.03	0.07	0.16	0.09	0.08	6	0.03	0.08	0.12	0.16	0.04
2,4,4-Trimethyl-1-Pentene	<DL	<DL	<DL	0.01	<DL	0.04	6	<DL	0.01	0.02	0.04	0.02
2,4-4-Trimethyl-2-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2,4-Dimethylpentane	0.07	0.04	0.06	0.17	0.09	0.06	6	0.04	0.08	0.12	0.17	0.05
2,5-Dimethylhexane	0.03	<DL	0.02	0.09	0.04	0.03	6	<DL	0.04	0.06	0.09	0.03
2-Ethyl-1-Butene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2-Methyl-1-Pentene	0.05	0.01	0.02	0.05	0.03	0.03	6	0.01	0.03	0.04	0.05	0.02
2-Methyl-2-Pentene	0.08	<DL	0.03	0.07	0.05	0.02	6	<DL	0.04	0.07	0.08	0.03
2-Methylheptane	0.07	0.04	0.07	0.1	0.1	0.1	6	0.04	0.08	0.10	0.10	0.02
2-Propanol	0.85	0.44	1.21	0.36	12.7	3.16	6	0.36	3.12	7.04	12.70	4.80
3-Methyl-1-Butene	0.07	<DL	<DL	0.08	0.07	0.04	6	<DL	0.04	0.07	0.08	0.04
3-Methylheptane	0.04	<DL	0.03	0.07	0.04	0.04	6	<DL	0.04	0.06	0.07	0.02
3-Methylhexane	0.24	0.15	0.26	0.4	0.24	0.33	6	0.15	0.27	0.34	0.40	0.09
3-Methylpentane	0.41	0.14	0.3	0.54	0.37	0.27	6	0.14	0.34	0.45	0.54	0.14
4-Methyl-1-Pentene	0.03	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.03	0.01
4-Nonene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Acetaldehyde	4.16	3.03	4.86	4.43	4.36	9.63	6	3.03	5.08	6.97	9.63	2.31
Acetone (+)	3.82	3.81	4.22	4.07	4.06	11.4	6	3.81	5.23	7.70	11.40	3.03
Acetonitrile	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Acetylene	0.98	0.54	0.74	0.91	0.89	0.64	6	0.54	0.78	0.92	0.98	0.17
Acrylonitrile	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Allyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
a-Pinene	0.07	0.06	0.07	0.06	0.06	0.06	6	0.06	0.06	0.07	0.07	0.01
Benzaldehyde	0.18	<DL	0.23	0.26	0.13	0.42	6	<DL	0.20	0.32	0.42	0.14
Benzene	0.47	0.2	0.31	0.63	0.46	0.37	6	0.20	0.41	0.53	0.63	0.15
Benzyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
b-Pinene	0.02	0.02	0.02	0.02	0.02	0.02	6	0.02	0.02	0.02	0.02	0.00
Bromochloromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromodichloromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromoform	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromomethane	0.03	<DL	<DL	<DL	0.02	0.02	6	<DL	0.01	0.02	0.03	0.01
Butyl Acrylate	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Butyraldehyde	1.51	0.49	0.46	0.36	0.51	0.83	6	0.36	0.69	1.04	1.51	0.43
c-1,2-Dichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
c-2-Butene	0.1	<DL	0.03	0.13	0.33	0.07	6	<DL	0.11	0.21	0.33	0.12
c-2-Hexene	0.03	<DL	<DL	0.02	0.02	<DL	6	<DL	0.01	0.02	0.03	0.01
c-2-Octene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC

Phase I Sampling Program - 10/23/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
c-2-Pentene	0.14	0.03	0.05	0.14	0.13	0.05	6	0.03	0.09	0.13	0.14	0.05
c-3-Hexene	0.06	<DL	<DL	0.03	0.04	<DL	6	<DL	0.02	0.04	0.06	0.03
c-3-Methyl-2-Pentene	0.05	<DL	<DL	0.03	<DL	<DL	6	<DL	0.01	0.03	0.05	0.02
c-4-Methyl-2-Pentene	0.04	<DL	<DL	0.03	0.02	<DL	6	<DL	0.02	0.03	0.04	0.02
Carbon Tetrachloride	0.11	0.12	0.11	0.11	0.12	0.1	6	0.10	0.11	0.12	0.12	0.01
Chlorobenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Chlorodifluoromethane	0.29	0.3	0.34	0.45	0.45	0.43	6	0.29	0.38	0.44	0.45	0.08
Chloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Chloroform	0.03	0.02	<DL	<DL	0.02	0.03	6	<DL	0.02	0.03	0.03	0.01
Chloromethane	0.79	0.87	0.84	0.81	0.88	0.97	6	0.79	0.86	0.91	0.97	0.06
Chloroprene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Cyclohexane	0.18	0.12	0.26	0.34	0.2	0.23	6	0.12	0.22	0.28	0.34	0.07
Cyclohexene	<DL	0.03	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.03	0.01
Cyclopentane	0.11	0.05	0.09	0.18	0.11	0.1	6	0.05	0.11	0.14	0.18	0.04
Cyclopentene	0.06	<DL	0.03	0.12	0.06	0.04	6	<DL	0.05	0.08	0.12	0.04
Dibromochloromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Dichlorodifluoromethane	0.62	0.64	0.66	0.63	0.65	0.64	6	0.62	0.64	0.65	0.66	0.01
Dichlorofluoromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Diethyl Ether	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Epichlorohydrin	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Ethane	53	5.86	7.66	8.88	8.58	6.2	6	5.86	15.03	30.25	53.00	18.64
Ethanol	4.83	2.65	6.42	2.52	5.38	11.9	6	2.52	5.62	8.43	11.90	3.44
Ethylbenzene	0.14	0.03	0.1	0.1	0.1	0.08	6	0.03	0.09	0.12	0.14	0.04
Ethylene	2.46	0.79	1.31	2.87	1.6	1.81	6	0.79	1.81	2.43	2.87	0.76
Freon 113	0.13	0.1	0.1	0.09	0.1	0.09	6	0.09	0.10	0.11	0.13	0.01
Freon 114	0.03	0.02	0.01	0.02	0.02	0.02	6	0.01	0.02	0.03	0.03	0.01
Halocarbon 134A	0.8	0.07	0.08	0.07	0.08	0.94	6	0.07	0.34	0.68	0.94	0.41
Heptanal	0.23	<DL	<DL	<DL	<DL	0.48	6	<DL	0.12	0.28	0.48	0.20
Hexachloro-1,3-Butadiene*	0.03	<DL	0.04	<DL	<DL	<DL	6	<DL	0.01	0.03	0.04	0.02
Hexanal	0.83	0.23	0.64	<DL	0.11	0.77	6	<DL	0.43	0.72	0.83	0.36
Indan	0.02	<DL	<DL	0.01	<DL	<DL	6	<DL	0.01	0.01	0.02	0.01
Indene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Isobutane	1.56	0.6	1.25	3.28	2.52	2.01	6	0.60	1.87	2.65	3.28	0.95
Isobutene + 1-Butene	0.34	0.2	0.21	0.32	0.84	0.33	6	0.20	0.37	0.57	0.84	0.24
Isobutylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Isoheptane	0.16	0.1	0.18	0.4	0.22	0.4	6	0.10	0.24	0.35	0.40	0.13
Isohexane	0.63	0.24	0.42	0.92	0.64	0.48	6	0.24	0.56	0.74	0.92	0.23
Isoprene	0.09	0.05	0.1	0.19	0.12	0.18	6	0.05	0.12	0.17	0.19	0.05
m-Dichlorobenzene	0.03	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.03	0.01
m-Diethylbenzene	0.01	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	0.01	NC
Methyl ethyl ketone	0.87	0.69	0.83	0.75	0.69	1.08	6	0.69	0.82	0.94	1.08	0.15
Methyl t-Butylether	0.13	0.21	0.13	0.32	0.19	0.23	6	0.13	0.20	0.26	0.32	0.07
Methylcyclohexane	0.1	0.08	0.17	0.41	0.23	0.35	6	0.08	0.22	0.33	0.41	0.13
Methylcyclopentane	0.28	0.14	0.29	0.65	0.34	0.33	6	0.14	0.34	0.48	0.65	0.17
Methylene Chloride	0.15	0.07	0.18	0.13	0.12	0.84	6	0.07	0.25	0.49	0.84	0.29
Methylisobutylketone	<DL	<DL	<DL	<DL	<DL	0.2	6	<DL	0.03	0.10	0.20	0.08
Naphthalene	<DL	<DL	<DL	<DL	<DL	0.05	6	<DL	0.01	0.03	0.05	0.02
n-Butane	3.43	1.3	2.2	4.98	4.83	2.82	6	1.30	3.26	4.45	4.98	1.46
n-Butylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
n-Decane	0.04	0.01	0.07	0.06	0.05	0.07	6	0.01	0.05	0.07	0.07	0.02
Neohexane	0.09	0.04	0.07	0.08	0.08	0.07	6	0.04	0.07	0.09	0.09	0.02
Neopentane	0.03	0.01	0.02	0.03	0.02	0.02	6	0.01	0.02	0.03	0.03	0.01
n-Heptane	0.12	0.06	0.15	0.27	0.17	0.23	6	0.06	0.17	0.23	0.27	0.08
n-Hexane	0.39	0.18	0.34	0.63	0.41	0.41	6	0.18	0.39	0.51	0.63	0.14
n-Nonane	0.06	0.05	0.06	0.08	0.08	0.06	6	0.05	0.07	0.08	0.08	0.01
n-Octane	0.08	0.05	0.12	0.14	0.11	0.12	6	0.05	0.10	0.13	0.14	0.03
n-Pentane	1.23	0.53	1.08	1.82	1.14	1.44	6	0.53	1.21	1.55	1.82	0.43
n-Propylbenzene	0.02	<DL	<DL	0.02	<DL	0.02	6	<DL	0.01	0.02	0.02	0.01
n-Undecane	0.05	<DL	0.07	0.05	0.05	0.09	6	<DL	0.05	0.08	0.09	0.03
o-Dichlorobenzene	0.03	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.03	0.01
p-Dichlorobenzene	0.14	0.05	0.21	0.12	0.1	1.02	6	0.05	0.27	0.58	1.02	0.37
p-Diethylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
p-Isopropyltoluene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Propane	5.25	2.68	4	11.6	7.06	5.86	6	2.68	6.08	8.61	11.60	3.10
Propylene	3.24	0.36	0.69	3.81	4.32	1.68	6	0.36	2.35	3.71	4.32	1.67
Styrene	0.08	0.02	0.08	0.03	0.03	0.08	6	0.02	0.05	0.08	0.08	0.03

Phase I Sampling Program - 10/23/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
t-1,2-Dichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
t-2-Butene	0.09	0.02	0.04	0.11	0.4	0.06	6	0.02	0.12	0.24	0.40	0.14
t-2-Hexene	0.06	<DL	<DL	0.05	0.04	0.02	6	<DL	0.03	0.05	0.06	0.03
t-2-Pentene	0.25	0.04	0.08	0.3	0.22	0.1	6	0.04	0.17	0.25	0.30	0.11
t-Butylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Tetrachloroethylene	0.01	<DL	<DL	<DL	<DL	0.04	6	<DL	0.01	0.02	0.04	0.02
Toluene	0.78	0.24	0.61	0.77	0.63	0.76	6	0.24	0.63	0.80	0.78	0.21
Trichloroethylene	0.24	0.03	0.29	0.18	0.31	2.94	6	0.03	0.67	1.58	2.94	1.12
Trichlorofluoromethane	0.37	0.31	0.33	0.29	0.31	0.38	6	0.29	0.33	0.36	0.38	0.04
Vinyl Acetate	<DL	0.71	<DL	<DL	<DL	<DL	6	<DL	0.12	0.36	0.71	0.29
Vinyl Bromide	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Vinyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,3-Dichloropropene (Total)	ND	ND	ND	ND	ND	ND	6	<DL	NC	NC	<DL	NC
Xylene (Total)	0.55	0.13	0.39	0.43	0.37	0.28	6	0.13	0.36	0.47	0.55	0.14

<DL- Less than Detection Limit

Min- Minimum Value

Max- Maximum Value

NC- Not Calculated

STD- Standard Deviation

* Suspected Laboratory Contaminant

Phase I Sampling Program - 10/29/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
1,1,1-Trichloroethane	<DL	0.03	0.03	0.03	0.03	0.03	6	<DL	0.03	0.04	0.03	0.01
1,1,2,2-Tetrachloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,1,2-Trichloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,1-Dichloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,1-Dichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2,3-Trimethylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2,4-Trichlorobenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2,4-Trimethylbenzene	<DL	<DL	0.03	0.07	0.04	0.02	6	<DL	0.03	0.05	0.07	0.03
1,2-Dibromoethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2-Dichloroethane	<DL	0.11	<DL	<DL	<DL	<DL	6	<DL	0.02	0.06	0.11	0.04
1,2-Dichloropropane	<DL	0.14	<DL	<DL	<DL	<DL	6	<DL	0.02	0.07	0.14	0.06
1,3,5-Trimethylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,3-Butadiene	0.55	0.24	0.2	0.3	0.2	0.21	6	0.20	0.28	0.39	0.55	0.14
1,4-Dioxane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Butanol	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Decene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Heptene	<DL	0.02	<DL	<DL	<DL	<DL	6	<DL	NC	NC	0.02	NC
1-Hexene	<DL	0.07	<DL	<DL	<DL	0.19	6	<DL	0.04	0.11	0.19	0.08
1-Methylcyclohexene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Nonene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Octene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Pentene	0.27	0.08	0.07	0.15	0.1	0.07	6	0.07	0.12	0.19	0.27	0.08
1-Propanol	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Undecene	<DL	0.09	0.08	0.07	<DL	0.06	6	<DL	0.05	0.08	0.09	0.04
2,2,3-Trimethylpentane	<DL	<DL	<DL	0.01	<DL	<DL	6	<DL	NC	NC	0.01	NC
2,2,4-Trimethylpentane	0.1	0.05	0.06	0.11	0.08	0.05	6	0.05	0.08	0.10	0.11	0.03
2,2,5-Trimethylhexane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2,3,4-Trimethylpentane	<DL	<DL	<DL	0.11	<DL	<DL	6	<DL	0.02	0.06	0.11	0.04
2,3-Dimethylbutane	0.15	0.06	0.07	0.09	0.07	0.06	6	0.06	0.08	0.11	0.15	0.03
2,3-Dimethylpentane	0.11	0.02	0.03	0.05	0.05	0.02	6	0.02	0.05	0.07	0.11	0.03
2,4,4-Trimethyl-1-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2,4-4-Trimethyl-2-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2,4-Dimethylpentane	<DL	<DL	<DL	0.04	0.03	<DL	6	<DL	0.01	0.03	0.04	0.02
2,5-Dimethylhexane	<DL	<DL	<DL	0.01	<DL	<DL	6	<DL	NC	NC	0.01	NC
2-Ethyl-1-Butene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2-Methyl-1-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2-Methyl-2-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2-Methylheptane	0.08	<DL	<DL	0.04	0.02	<DL	6	<DL	0.02	0.05	0.08	0.03
2-Propanol	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
3-Methyl-1-Butene	<DL	<DL	<DL	<DL	0.04	<DL	6	<DL	0.01	0.02	0.04	0.02
3-Methylheptane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
3-Methylhexane	0.16	0.06	0.07	0.11	0.11	0.09	6	0.06	0.10	0.13	0.16	0.04
3-Methylpentane	0.29	0.1	0.11	0.87	0.71	0.68	6	0.10	0.46	0.73	0.87	0.33
4-Methyl-1-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
4-Nonene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Acetaldehyde	12.5	2.19	2.56	4.62	1.75	6.7	6	1.75	5.05	8.39	12.50	4.09
Acetone (+)	6.71	1.56	1.61	1	1.21	1.54	6	1.00	2.27	4.06	6.71	2.19
Acetonitrile	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Acetylene	1.42	0.71	0.78	1.16	0.99	0.85	6	0.71	0.99	1.20	1.42	0.27
Acrylonitrile	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Allyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
a-Pinene	<DL	0.04	0.03	0.03	0.02	0.03	6	<DL	0.03	0.04	0.04	0.01
Benzaldehyde	<DL	<DL	<DL	0.13	<DL	<DL	6	<DL	0.02	0.07	0.13	0.05
Benzene	0.45	0.32	0.24	0.49	0.31	0.26	6	0.24	0.35	0.43	0.49	0.10
Benzyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
b-Pinene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromochloromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromodichloromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromoform	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromomethane	0.11	<DL	<DL	<DL	<DL	<DL	6	<DL	0.02	0.06	0.11	0.04
Butyl Acrylate	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Butyraldehyde	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
c-1,2-Dichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
c-2-Butene	0.41	0.23	0.09	0.21	0.21	0.12	6	0.09	0.21	0.30	0.41	0.11
c-2-Hexene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
c-2-Octene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC

Phase I Sampling Program - 10/29/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
c-2-Pentene	0.06	<DL	0.01	0.05	0.04	<DL	6	<DL	0.03	0.05	0.06	0.03
c-3-Hexene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
c-3-Methyl-2-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
c-4-Methyl-2-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Carbon Tetrachloride	0.08	0.11	0.11	0.11	0.12	0.12	6	0.08	0.11	0.12	0.12	0.01
Chlorobenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Chlorodifluoromethane	0.6	0.29	0.56	0.62	0.42	0.45	6	0.29	0.49	0.59	0.62	0.13
Chloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Chloroform	<DL	<DL	<DL	0.02	0.04	0.02	6	<DL	0.01	0.03	0.04	0.02
Chloromethane	0.98	0.73	0.79	0.85	0.85	0.74	6	0.73	0.82	0.90	0.98	0.09
Chloroprene	0.13	0.24	0.16	0.4	0.47	0.56	6	0.13	0.33	0.47	0.56	0.18
Cyclohexane	<DL	0.04	0.03	0.07	0.04	0.03	6	<DL	0.04	0.05	0.07	0.02
Cyclohexene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Cyclopentane	<DL	<DL	<DL	0.05	0.04	<DL	6	<DL	0.02	0.03	0.05	0.02
Cyclopentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Dibromochloromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Dichlorodifluoromethane	0.78	0.64	0.69	0.71	0.72	0.73	6	0.64	0.71	0.75	0.78	0.05
Dichlorofluoromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Diethyl Ether	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Epichlorohydrin	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Ethane	16	5.34	6.48	6.99	9.79	9.04	6	5.34	8.94	12.07	16.00	3.83
Ethanol	5.37	<DL	<DL	<DL	<DL	<DL	6	<DL	0.90	2.69	5.37	2.19
Ethylbenzene	<DL	<DL	0.04	0.07	0.05	<DL	6	<DL	0.03	0.05	0.07	0.03
Ethylene	4.88	3.3	2.51	4.19	2.53	2.45	6	2.45	3.31	4.14	4.88	1.02
Freon 113	0.17	0.1	0.1	0.1	0.09	0.09	6	0.09	0.11	0.13	0.17	0.03
Freon 114	0.07	<DL	<DL	0.01	<DL	<DL	6	<DL	0.01	0.04	0.07	0.03
Halocarbon 134A	0.1	0.03	0.12	0.13	0.19	0.07	6	0.03	0.11	0.15	0.19	0.05
Heptanal	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Hexachloro-1,3-Butadiene*	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Hexanal	<DL	0.2	<DL	0.1	<DL	<DL	6	<DL	0.05	0.12	0.20	0.08
Indan	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Indene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Isobutane	2.92	1.1	1.3	1.98	1.8	1.32	6	1.10	1.74	2.28	2.92	0.67
Isobutene + 1-Butene	1.6	0.82	0.26	0.6	0.42	0.54	6	0.26	0.71	1.10	1.60	0.48
Isobutylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Isoheptane	0.21	0.06	0.07	0.11	0.12	0.08	6	0.06	0.11	0.15	0.21	0.05
Isohexane	0.47	0.19	0.24	0.45	0.34	0.22	6	0.19	0.32	0.42	0.47	0.12
Isoprene	<DL	<DL	0.02	0.13	0.03	<DL	6	<DL	0.03	0.07	0.13	0.05
m-Dichlorobenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
m-Diethylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Methyl ethyl ketone	<DL	2.42	5.34	10.3	18.8	20.3	6	<DL	9.53	16.47	20.30	8.50
Methyl t-Butylether	<DL	<DL	<DL	0.02	<DL	<DL	6	<DL	NC	NC	0.02	NC
Methylcyclohexane	0.04	0.04	0.02	0.06	0.04	0.03	6	0.02	0.04	0.05	0.06	0.01
Methylcyclopentane	0.15	0.07	0.08	0.14	0.1	0.06	6	0.06	0.10	0.13	0.15	0.04
Methylene Chloride	1.65	0.14	0.31	0.19	0.21	0.15	6	0.14	0.44	0.93	1.65	0.60
Methylisobutylketone	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Naphthalene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
n-Butane	5.8	2.18	2.27	4.45	3.58	2.72	6	2.18	3.50	4.66	5.80	1.42
n-Butylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
n-Decane	0.17	0.01	0.01	0.02	0.01	<DL	6	<DL	0.04	0.09	0.17	0.07
Neohexane	0.08	0.03	0.12	0.05	0.11	0.04	6	0.03	0.07	0.10	0.12	0.04
Neopentane	<DL	0.01	0.01	0.02	0.01	0.01	6	<DL	0.01	0.02	0.02	0.01
n-Heptane	0.23	0.03	0.04	0.07	0.05	0.02	6	0.02	0.07	0.14	0.23	0.08
n-Hexane	0.55	0.18	0.24	0.39	0.28	0.21	6	0.18	0.31	0.42	0.55	0.14
n-Nonane	0.34	<DL	0.02	0.03	<DL	<DL	6	<DL	0.07	0.18	0.34	0.14
n-Octane	0.28	0.03	0.03	0.05	0.03	0.02	6	0.02	0.07	0.16	0.28	0.10
n-Pentane	2.77	0.58	0.89	1.22	1.02	0.95	6	0.58	1.24	1.87	2.77	0.78
n-Propylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
n-Undecane	0.12	0.02	0.01	0.02	0.02	<DL	6	<DL	0.03	0.07	0.12	0.04
o-Dichlorobenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
p-Dichlorobenzene	1.03	0.23	0.16	0.07	0.07	0.08	6	0.07	0.27	0.58	1.03	0.38
p-Diethylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
p-Isopropyltoluene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Propane	9.42	4.58	4.72	4.83	4.12	4	6	4.00	5.28	6.96	9.42	2.06
Propylene	7.03	2.11	2.09	3.58	2.37	1.74	6	1.74	3.15	4.79	7.03	2.00
Styrene	<DL	0.05	0.05	0.04	0.03	<DL	6	<DL	0.03	0.05	0.05	0.02

Phase I Sampling Program - 10/29/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
t-1,2-Dichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
t-2-Butene	0.49	0.33	0.17	0.21	0.31	0.18	6	0.17	0.28	0.38	0.49	0.12
t-2-Hexene	<DL	<DL	<DL	0.02	<DL	<DL	6	<DL	NC	NC	0.02	NC
t-2-Pentene	0.17	0.02	0.05	0.11	0.07	0.03	6	0.02	0.08	0.12	0.17	0.06
t-Butylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Tetrachloroethylene	<DL	<DL	<DL	0.01	<DL	<DL	6	<DL	NC	NC	0.01	NC
Toluene	0.98	0.82	0.81	1.24	0.94	0.79	6	0.79	0.93	1.07	1.24	0.17
Trichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Trichlorofluoromethane	0.51	0.33	0.33	0.33	0.33	0.33	6	0.33	0.36	0.42	0.51	0.07
Vinyl Acetate	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Vinyl Bromide	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Vinyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,3-Dichloropropene (Total)	ND	0.09	ND	ND	ND	ND	6	<DL	0.02	0.05	0.09	0.04
Xylene (Total)	0.32	0.12	0.18	0.31	0.23	0.16	6	0.12	0.22	0.29	0.32	0.08

<DL- Less than Detection Limit

Min- Minimum Value

Max- Maximum Value

NC- Not Calculated

STD- Standard Deviation

* Suspected Laboratory Contaminant

Phase I Sampling Program - 11/04/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
1,1,1-Trichloroethane	0.03	0.03	0.05	0.06	0.03	0.03	6	0.03	0.04	0.05	0.06	0.01
1,1,2,2-Tetrachloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,1,2-Trichloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,1-Dichloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,1-Dichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2,3-Trimethylbenzene	0.03	<DL	0.02	0.03	0.02	0.02	6	<DL	0.02	0.03	0.03	0.01
1,2,4-Trichlorobenzene	<DL	<DL	0.23	0.01	<DL	<DL	6	<DL	0.04	0.12	0.23	0.09
1,2,4-Trimethylbenzene	0.11	0.07	0.09	0.12	0.09	0.07	6	0.07	0.09	0.11	0.12	0.02
1,2-Dibromoethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2-Dichloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2-Dichloropropane	<DL	0.03	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.03	0.01
1,3,5-Trimethylbenzene	0.04	0.02	0.02	0.04	0.03	0.02	6	0.02	0.03	0.04	0.04	0.01
1,3-Butadiene	0.31	0.15	0.13	0.28	0.32	0.17	6	0.13	0.23	0.30	0.32	0.09
1,4-Dioxane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Butanol	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Decene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Heptene	0.21	0.07	0.13	0.07	0.08	0.07	6	0.07	0.11	0.15	0.21	0.06
1-Hexene	0.05	0.05	0.07	0.07	0.03	0.26	6	0.03	0.09	0.16	0.26	0.09
1-Methylcyclohexene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Nonene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Octene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Pentene	0.25	0.13	0.17	0.24	0.16	0.18	6	0.13	0.19	0.23	0.25	0.05
1-Propanol	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Undecene	<DL	0.06	0.05	0.2	<DL	0.04	6	<DL	0.06	0.12	0.20	0.07
2,2,3-Trimethylpentane	0.03	0.02	0.03	0.06	0.02	0.02	6	0.02	0.03	0.04	0.06	0.02
2,2,4-Trimethylpentane	0.23	0.16	0.22	0.4	0.19	0.2	6	0.16	0.23	0.30	0.40	0.09
2,2,5-Trimethylhexane	0.02	0.01	0.02	0.04	0.02	0.01	6	0.01	0.02	0.03	0.04	0.01
2,3,4-Trimethylpentane	0.15	0.13	0.15	0.3	0.14	0.14	6	0.13	0.17	0.22	0.30	0.06
2,3-Dimethylbutane	0.23	0.12	0.15	0.19	0.15	0.15	6	0.12	0.17	0.20	0.23	0.04
2,3-Dimethylpentane	0.11	0.06	0.09	0.11	0.08	0.07	6	0.06	0.09	0.10	0.11	0.02
2,4,4-Trimethyl-1-Pentene	<DL	<DL	<DL	0.02	<DL	<DL	6	<DL	NC	NC	0.02	NC
2,4-4-Trimethyl-2-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2,4-Dimethylpentane	0.07	0.04	0.05	0.06	0.06	0.05	6	0.04	0.06	0.06	0.07	0.01
2,5-Dimethylhexane	0.03	0.02	0.03	0.07	0.03	0.03	6	0.02	0.04	0.05	0.07	0.02
2-Ethyl-1-Butene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2-Methyl-1-Pentene	0.03	0.03	0.04	0.04	0.03	<DL	6	<DL	0.03	0.04	0.04	0.01
2-Methyl-2-Pentene	0.05	0.03	0.04	0.05	0.03	0.05	6	0.03	0.04	0.05	0.05	0.01
2-Methylheptane	0.09	0.06	0.08	0.12	0.07	0.07	6	0.06	0.08	0.10	0.12	0.02
2-Propanol	5.8	1.58	0.41	0.45	<DL	0.99	6	<DL	1.54	3.30	5.80	2.16
3-Methyl-1-Butene	0.1	0.04	0.06	0.09	0.06	0.07	6	0.04	0.07	0.09	0.10	0.02
3-Methylheptane	0.05	0.03	0.04	0.08	0.05	0.03	6	0.03	0.05	0.06	0.08	0.02
3-Methylhexane	0.3	0.2	0.25	0.34	0.19	0.22	6	0.19	0.25	0.30	0.34	0.06
3-Methylpentane	0.55	0.32	0.39	0.43	0.33	0.4	6	0.32	0.40	0.47	0.55	0.08
4-Methyl-1-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
4-Nonene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Acetaldehyde	2.03	6.71	6.38	5.36	3.09	5.42	6	2.03	4.83	6.36	6.71	1.87
Acetone (+)	3.28	3.84	5.07	5.54	2.71	2.82	6	2.71	3.88	4.84	5.54	1.19
Acetonitrile	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Acetylene	0.99	0.82	0.83	0.89	1.12	0.82	6	0.82	0.91	1.01	1.12	0.12
Acrylonitrile	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Allyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
a-Pinene	0.07	0.07	0.07	0.08	0.08	0.07	6	0.07	0.07	0.08	0.08	0.01
Benzaldehyde	<DL	0.27	0.3	0.42	0.18	<DL	6	<DL	0.20	0.33	0.42	0.17
Benzene	0.91	0.49	0.58	0.81	0.77	0.54	6	0.49	0.68	0.82	0.91	0.17
Benzyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
b-Pinene	0.04	0.03	0.05	0.04	0.05	0.04	6	0.03	0.04	0.05	0.05	0.01
Bromochloromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromodichloromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromoform	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromomethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Butyl Acrylate	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Butyraldehyde	0.18	0.37	0.44	0.42	0.18	0.15	6	0.15	0.29	0.40	0.44	0.13
c-1,2-Dichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
c-2-Butene	0.37	0.85	0.24	0.32	0.3	0.48	6	0.24	0.43	0.61	0.85	0.22
c-2-Hexene	0.02	<DL	0.02	0.02	<DL	<DL	6	<DL	0.01	0.02	0.02	0.01
c-2-Octene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC

Phase I Sampling Program - 11/04/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
c-2-Pentene	0.18	0.09	0.12	0.15	0.1	0.14	6	0.09	0.13	0.16	0.18	0.03
c-3-Hexene	0.03	<DL	0.01	0.02	<DL	<DL	6	<DL	0.01	0.02	0.03	0.01
c-3-Methyl-2-Pentene	0.01	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	0.01	NC
c-4-Methyl-2-Pentene	0.03	<DL	0.01	<DL	<DL	0.02	6	<DL	0.01	0.02	0.03	0.01
Carbon Tetrachloride	0.11	0.12	0.12	0.1	0.12	0.12	6	0.10	0.12	0.12	0.12	0.01
Chlorobenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Chlorodifluoromethane	0.3	0.3	0.39	0.51	0.29	0.32	6	0.29	0.35	0.42	0.51	0.09
Chloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Chloroform	0.02	0.04	0.02	0.03	0.03	0.02	6	0.02	0.03	0.03	0.04	0.01
Chloromethane	0.92	0.91	1	0.77	0.7	0.78	6	0.70	0.85	0.94	1.00	0.11
Chloroprene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Cyclohexane	0.21	0.12	0.15	0.18	0.12	0.14	6	0.12	0.15	0.18	0.21	0.04
Cyclohexene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Cyclopentane	0.17	0.08	0.1	0.12	0.09	0.1	6	0.08	0.11	0.14	0.17	0.03
Cyclopentene	0.17	0.05	0.08	0.1	0.05	0.06	6	0.05	0.09	0.12	0.17	0.05
Dibromochloromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Dichlorodifluoromethane	0.65	0.7	0.66	0.67	0.67	0.68	6	0.65	0.67	0.69	0.70	0.02
Dichlorofluoromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Diethyl Ether	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Epichlorohydrin	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Ethane	64.7	7.96	8.26	8.59	8.94	7.15	6	7.15	17.60	36.45	64.70	23.08
Ethanol	0.95	1.08	1.74	1.66	0.79	1.07	6	0.79	1.22	1.53	1.74	0.39
Ethylbenzene	0.13	0.08	0.11	0.15	0.1	0.09	6	0.08	0.11	0.13	0.15	0.03
Ethylene	8.17	2.42	2.2	2.62	3.08	3.06	6	2.20	3.59	5.44	8.17	2.27
Freon 113	0.09	0.1	0.09	0.09	0.1	0.1	6	0.09	0.10	0.10	0.10	0.01
Freon 114	0.01	<DL	0.02	0.02	0.01	0.01	6	<DL	0.01	0.02	0.02	0.01
Halocarbon 134A	0.43	0.13	4.12	6.72	0.11	0.82	6	0.11	2.06	4.30	6.72	2.75
Heptanal	<DL	0.14	<DL	0.21	<DL	<DL	6	<DL	0.06	0.13	0.21	0.09
Hexachloro-1,3-Butadiene*	<DL	<DL	0.21	<DL	<DL	<DL	6	<DL	0.04	0.11	0.21	0.09
Hexanal	0.15	0.29	0.3	0.61	0.17	0.12	6	0.12	0.27	0.42	0.61	0.18
Indan	<DL	<DL	<DL	0.02	<DL	<DL	6	<DL	NC	NC	0.02	NC
Indene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Isobutane	4.71	2.05	2.65	5.39	2.75	3.13	6	2.05	3.45	4.51	5.39	1.31
Isobutene + 1-Butene	1.08	0.85	0.84	1.02	1	1.3	6	0.84	1.02	1.15	1.30	0.17
Isobutylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Isoheptane	0.25	0.14	0.18	0.24	0.15	0.18	6	0.14	0.19	0.23	0.25	0.05
Isohexane	0.89	0.52	0.59	0.71	0.53	0.61	6	0.52	0.64	0.76	0.89	0.14
Isoprene	0.22	0.08	0.13	0.3	0.14	0.1	6	0.08	0.16	0.23	0.30	0.08
m-Dichlorobenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
m-Diethylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Methyl ethyl ketone	0.2	1.92	0.66	0.72	0.49	0.22	6	0.20	0.70	1.22	1.92	0.63
Methyl t-Butylether	0.23	0.09	0.1	0.08	0.07	0.07	6	0.07	0.11	0.16	0.23	0.06
Methylcyclohexane	0.28	0.16	0.26	0.39	0.17	0.17	6	0.16	0.24	0.31	0.39	0.09
Methylcyclopentane	0.42	0.21	0.24	0.32	0.22	0.24	6	0.21	0.28	0.34	0.42	0.08
Methylene Chloride	0.22	0.11	0.17	0.4	0.09	0.13	6	0.09	0.19	0.28	0.40	0.11
Methylisobutylketone	<DL	0.41	0.05	0.07	<DL	<DL	6	<DL	0.09	0.22	0.41	0.16
Naphthalene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
n-Butane	8.38	4.78	5.24	6.97	5.01	5.76	6	4.78	6.02	7.16	8.38	1.39
n-Butylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
n-Decane	0.05	0.04	0.08	0.16	0.05	0.05	6	0.04	0.07	0.11	0.16	0.05
Neohexane	0.1	0.08	0.07	0.08	0.06	0.07	6	0.06	0.08	0.09	0.10	0.01
Neopentane	0.06	0.02	0.03	0.04	0.03	0.03	6	0.02	0.04	0.05	0.06	0.01
n-Heptane	0.24	0.16	0.23	0.37	0.18	0.22	6	0.16	0.23	0.29	0.37	0.07
n-Hexane	0.65	0.46	0.5	0.66	0.42	0.5	6	0.42	0.53	0.61	0.66	0.10
n-Nonane	0.08	0.04	0.07	0.14	0.07	0.06	6	0.04	0.08	0.10	0.14	0.03
n-Octane	0.12	0.09	0.13	0.25	0.13	0.1	6	0.09	0.14	0.18	0.25	0.06
n-Pentane	2.73	1.28	1.47	1.98	1.38	1.48	6	1.28	1.72	2.17	2.73	0.55
n-Propylbenzene	<DL	<DL	<DL	0.02	<DL	<DL	6	<DL	NC	NC	0.02	NC
n-Undecane	0.03	0.03	0.05	0.11	0.03	0.03	6	0.03	0.05	0.07	0.11	0.03
o-Dichlorobenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
p-Dichlorobenzene	0.16	0.1	0.84	1.39	0.13	0.2	6	0.10	0.47	0.90	1.39	0.53
p-Diethylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
p-Isopropyltoluene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Propane	11.7	7.39	9.36	14	11.4	7.58	6	7.39	10.24	12.35	14.00	2.59
Propylene	8.24	6.07	5.5	8.77	10.8	4.97	6	4.97	7.39	9.23	10.80	2.26
Styrene	0.06	0.04	0.04	0.36	0.04	0.04	6	0.04	0.10	0.20	0.36	0.13

Phase I Sampling Program - 11/04/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
t-1,2-Dichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
t-2-Butene	0.43	1.03	0.27	0.39	0.4	0.61	6	0.27	0.52	0.74	1.03	0.27
t-2-Hexene	0.04	0.03	0.04	0.05	0.03	0.04	6	0.03	0.04	0.04	0.05	0.01
t-2-Pentene	0.36	0.15	0.21	0.3	0.18	0.25	6	0.15	0.24	0.31	0.36	0.08
t-Butylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Tetrachloroethylene	<DL	<DL	0.02	0.02	0.02	<DL	6	<DL	0.01	0.02	0.02	0.01
Toluene	0.92	0.75	0.79	1.09	0.8	0.68	6	0.68	0.84	0.96	1.09	0.15
Trichloroethylene	<DL	<DL	0.05	0.19	0.03	<DL	6	<DL	0.05	0.11	0.19	0.07
Trichlorofluoromethane	0.34	0.34	0.61	0.79	0.32	0.38	6	0.32	0.46	0.62	0.79	0.19
Vinyl Acetate	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Vinyl Bromide	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Vinyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,3-Dichloropropene (Total)	ND	ND	ND	ND	ND	ND	6	<DL	NC	NC	<DL	NC
Xylene (Total)	0.54	0.33	0.4	0.53	0.42	0.36	6	0.33	0.43	0.50	0.54	0.09

<DL- Less than Detection Limit

Min- Minimum Value

Max- Maximum Value

NC- Not Calculated

STD- Standard Deviation

* Suspected Laboratory Contaminant

Phase I Sampling Program - 11/10/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
1,1,1-Trichloroethane	0.09	0.03	0.04	<DL	0.03	<DL	6	<DL	0.03	0.06	0.09	0.03
1,1,2,2-Tetrachloroethane	0.06	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.03	0.06	0.02
1,1,2-Trichloroethane	0.04	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.04	0.02
1,1-Dichloroethane	0.04	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.04	0.02
1,1-Dichloroethylene	0.05	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.03	0.05	0.02
1,2,3-Trimethylbenzene	0.06	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.03	0.06	0.02
1,2,4-Trichlorobenzene	0.15	0.03	<DL	0.02	<DL	<DL	6	<DL	0.03	0.08	0.15	0.06
1,2,4-Trimethylbenzene	0.09	0.03	0.02	0.08	0.04	0.02	6	0.02	0.05	0.07	0.09	0.03
1,2-Dibromoethane	0.05	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.03	0.05	0.02
1,2-Dichloroethane	0.06	0.19	<DL	<DL	<DL	<DL	6	<DL	0.04	0.10	0.19	0.08
1,2-Dichloropropane	0.04	0.27	<DL	<DL	<DL	<DL	6	<DL	0.05	0.14	0.27	0.11
1,3,5-Trimethylbenzene	0.06	0.01	<DL	0.02	<DL	<DL	6	<DL	0.02	0.03	0.06	0.02
1,3-Butadiene	0.19	0.34	0.08	0.05	0.04	0.04	6	0.04	0.12	0.22	0.34	0.12
1,4-Dioxane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Butanol	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Decene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Heptene	<DL	0.06	<DL	0.07	0.04	<DL	6	<DL	0.03	0.05	0.07	0.03
1-Hexene	0.18	0.72	0.23	0.1	0.08	0.07	6	0.07	0.23	0.43	0.72	0.25
1-Methylcyclohexene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Nonene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Octene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Pentene	0.1	0.07	0.05	0.08	0.07	0.04	6	0.04	0.07	0.09	0.10	0.02
1-Propanol	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Undecene	0.15	0.26	0.08	0.06	0.03	<DL	6	<DL	0.10	0.17	0.26	0.09
2,2,3-Trimethylpentane	<DL	<DL	<DL	0.01	<DL	<DL	6	<DL	NC	NC	0.01	NC
2,2,4-Trimethylpentane	0.12	0.05	0.04	0.12	0.07	0.04	6	0.04	0.07	0.10	0.12	0.04
2,2,5-Trimethylhexane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2,3,4-Trimethylpentane	0.12	<DL	<DL	<DL	0.08	<DL	6	<DL	0.03	0.08	0.12	0.05
2,3-Dimethylbutane	0.11	0.03	0.03	0.11	0.07	0.04	6	0.03	0.07	0.10	0.11	0.04
2,3-Dimethylpentane	0.09	0.06	0.02	0.05	0.03	0.02	6	0.02	0.05	0.07	0.09	0.03
2,4,4-Trimethyl-1-Pentene	0.02	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	0.02	NC
2,4-4-Trimethyl-2-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2,4-Dimethylpentane	0.07	<DL	<DL	0.03	0.02	<DL	6	<DL	0.02	0.04	0.07	0.03
2,5-Dimethylhexane	<DL	<DL	<DL	0.01	<DL	<DL	6	<DL	NC	NC	0.01	NC
2-Ethyl-1-Butene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2-Methyl-1-Pentene	0.05	0.41	0.13	0.01	<DL	0.05	6	<DL	0.11	0.23	0.41	0.15
2-Methyl-2-Pentene	<DL	0.02	<DL	0.02	0.01	<DL	6	<DL	0.01	0.02	0.02	0.01
2-Methylheptane	0.08	0.05	0.01	0.05	0.05	0.02	6	0.01	0.04	0.06	0.08	0.03
2-Propanol	1.24	1.04	1.03	0.23	3.18	0.15	6	0.15	1.15	2.04	3.18	1.10
3-Methyl-1-Butene	0.09	<DL	<DL	<DL	<DL	<DL	6	<DL	0.02	0.05	0.09	0.04
3-Methylheptane	0.04	<DL	<DL	<DL	0.01	<DL	6	<DL	0.01	0.02	0.04	0.02
3-Methylhexane	0.18	0.2	0.12	0.14	0.13	0.11	6	0.11	0.15	0.18	0.20	0.04
3-Methylpentane	0.15	0.09	0.09	0.22	0.17	0.1	6	0.09	0.14	0.18	0.22	0.05
4-Methyl-1-Pentene	0.05	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.03	0.05	0.02
4-Nonene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Acetaldehyde	5.03	5.56	4.84	3.09	4.24	2.69	6	2.69	4.24	5.17	5.56	1.14
Acetone (+)	2.92	3.29	2.69	2.22	3.66	2.12	6	2.12	2.82	3.31	3.66	0.60
Acetonitrile	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Acetylene	0.48	1.61	0.48	0.54	0.38	0.42	6	0.38	0.65	1.04	1.61	0.47
Acrylonitrile	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Allyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
a-Pinene	0.1	0.06	0.05	0.05	0.04	0.05	6	0.04	0.06	0.08	0.10	0.02
Benzaldehyde	0.28	0.12	0.2	0.09	0.13	0.28	6	0.09	0.18	0.25	0.28	0.08
Benzene	0.3	0.51	0.18	0.3	0.21	0.16	6	0.16	0.28	0.38	0.51	0.13
Benzyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
b-Pinene	0.05	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.03	0.05	0.02
Bromochloromethane	0.03	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.03	0.01
Bromodichloromethane	0.04	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.04	0.02
Bromoform	0.05	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.03	0.05	0.02
Bromomethane	0.08	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.04	0.08	0.03
Butyl Acrylate	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Butyraldehyde	0.39	0.22	0.28	0.18	0.34	0.38	6	0.18	0.30	0.37	0.39	0.09
c-1,2-Dichloroethylene	0.04	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.04	0.02
c-2-Butene	0.17	1.11	0.11	0.17	0.14	0.09	6	0.09	0.30	0.62	1.11	0.40
c-2-Hexene	0.03	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.03	0.01
c-2-Octene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC

Phase I Sampling Program - 11/10/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
c-2-Pentene	0.07	<DL	<DL	0.09	0.05	<DL	6	<DL	0.04	0.07	0.09	0.04
c-3-Hexene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
c-3-Methyl-2-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
c-4-Methyl-2-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Carbon Tetrachloride	0.19	0.12	0.12	0.12	0.13	0.12	6	0.12	0.13	0.16	0.19	0.03
Chlorobenzene	0.06	0.03	<DL	<DL	<DL	<DL	6	<DL	0.02	0.04	0.06	0.03
Chlorodifluoromethane	0.34	0.24	0.24	0.26	0.24	0.21	6	0.21	0.26	0.29	0.34	0.04
Chloroethane	0.19	<DL	<DL	<DL	<DL	<DL	6	<DL	0.03	0.10	0.19	0.08
Chloroform	0.05	0.05	0.01	<DL	0.01	<DL	6	<DL	0.02	0.04	0.05	0.02
Chloromethane	1.25	1.03	1	0.96	1.07	0.98	6	0.96	1.05	1.13	1.25	0.11
Chloroprene	0.04	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.04	0.02
Cyclohexane	0.11	0.18	0.04	0.09	0.08	0.05	6	0.04	0.09	0.13	0.18	0.05
Cyclohexene	<DL	0.02	<DL	<DL	<DL	<DL	6	<DL	NC	NC	0.02	NC
Cyclopentane	0.09	0.05	<DL	0.06	0.04	<DL	6	<DL	0.04	0.07	0.09	0.04
Cyclopentene	0.05	<DL	<DL	0.02	0.02	<DL	6	<DL	0.02	0.03	0.05	0.02
Dibromochloromethane	0.04	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.04	0.02
Dichlorodifluoromethane	0.76	0.68	0.69	0.7	0.69	0.67	6	0.67	0.70	0.72	0.76	0.03
Dichlorofluoromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Diethyl Ether	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Epichlorohydrin	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Ethane	4.18	5.19	3.74	3.19	3	3.75	6	3.00	3.84	4.48	5.19	0.79
Ethanol	1.39	<DL	<DL	<DL	0.56	<DL	6	<DL	0.33	0.79	1.39	0.57
Ethylbenzene	0.09	0.06	0.04	0.08	0.04	0.03	6	0.03	0.06	0.08	0.09	0.02
Ethylene	1.49	7.09	1.96	1.22	1.02	1.29	6	1.02	2.35	4.26	7.09	2.35
Freon 113	0.19	0.1	0.1	0.09	0.09	0.09	6	0.09	0.11	0.14	0.19	0.04
Freon 114	0.08	0.02	0.02	<DL	<DL	0.01	6	<DL	0.02	0.05	0.08	0.03
Halocarbon 134A	0.17	0.1	0.09	0.07	0.12	0.09	6	0.07	0.11	0.14	0.17	0.04
Heptanal	0.15	0.14	0.1	<DL	<DL	0.1	6	<DL	0.08	0.14	0.15	0.07
Hexachloro-1,3-Butadiene*	0.1	0.02	<DL	<DL	<DL	<DL	6	<DL	0.02	0.05	0.10	0.04
Hexanal	0.38	0.26	0.24	0.27	0.23	0.28	6	0.23	0.28	0.32	0.38	0.05
Indan	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Indene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Isobutane	0.83	1.05	0.87	1.11	0.99	0.8	6	0.80	0.94	1.05	1.11	0.13
Isobutene + 1-Butene	0.59	2.58	0.76	0.5	0.45	0.57	6	0.45	0.91	1.58	2.58	0.83
Isobutylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Isoheptane	0.1	0.11	0.06	0.1	0.08	0.05	6	0.05	0.08	0.10	0.11	0.02
Isohexane	0.23	0.12	0.14	0.36	0.25	0.14	6	0.12	0.21	0.28	0.36	0.09
Isoprene	0.16	0.05	0.09	0.11	0.06	0.04	6	0.04	0.09	0.12	0.16	0.05
m-Dichlorobenzene	0.07	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.04	0.07	0.03
m-Diethylbenzene	0.04	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.04	0.02
Methyl ethyl ketone	0.51	2.62	0.86	0.37	0.42	0.61	6	0.37	0.90	1.60	2.62	0.86
Methyl t-Butylether	0.51	<DL	<DL	<DL	0.04	<DL	6	<DL	0.09	0.26	0.51	0.21
Methylcyclohexane	0.11	0.23	0.08	0.13	0.13	0.06	6	0.06	0.12	0.17	0.23	0.06
Methylcyclopentane	0.13	0.1	0.06	0.15	0.1	0.06	6	0.06	0.10	0.13	0.15	0.04
Methylene Chloride	0.38	0.11	0.1	0.1	0.13	0.09	6	0.09	0.15	0.24	0.38	0.11
Methylisobutylketone	<DL	5.35	<DL	<DL	<DL	<DL	6	<DL	0.89	2.68	5.35	2.18
Naphthalene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
n-Butane	2.33	2.4	1.84	6.64	4.66	2.05	6	1.84	3.32	4.89	6.64	1.92
n-Butylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
n-Decane	0.06	0.05	0.02	0.02	0.03	<DL	6	<DL	0.03	0.05	0.06	0.02
Neohexane	0.08	0.02	<DL	0.05	0.04	0.02	6	<DL	0.04	0.06	0.08	0.03
Neopentane	<DL	<DL	<DL	0.02	0.01	<DL	6	<DL	0.01	0.01	0.02	0.01
n-Heptane	0.1	0.21	0.11	0.12	0.11	0.08	6	0.08	0.12	0.16	0.21	0.05
n-Hexane	0.25	0.28	0.17	0.27	0.22	0.13	6	0.13	0.22	0.27	0.28	0.06
n-Nonane	0.07	0.05	0.02	0.03	0.03	0.02	6	0.02	0.04	0.05	0.07	0.02
n-Octane	0.09	0.07	0.04	0.07	0.05	0.03	6	0.03	0.06	0.08	0.09	0.02
n-Pentane	0.52	0.57	0.37	1.02	1.12	0.38	6	0.37	0.66	0.93	1.12	0.33
n-Propylbenzene	0.05	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.03	0.05	0.02
n-Undecane	0.02	0.04	0.03	0.02	0.03	0.02	6	0.02	0.03	0.03	0.04	0.01
o-Dichlorobenzene	0.07	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.04	0.07	0.03
p-Dichlorobenzene	0.23	0.17	0.38	0.2	0.31	0.12	6	0.12	0.24	0.31	0.38	0.10
p-Diethylbenzene	0.05	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.03	0.05	0.02
p-Isopropyltoluene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Propane	2.12	2.83	2.15	2.09	2.47	2	6	2.00	2.28	2.53	2.83	0.31
Propylene	0.83	11.2	2.11	0.82	0.85	1.34	6	0.82	2.86	6.22	11.20	4.12
Styrene	0.09	0.17	0.06	0.04	<DL	0.18	6	<DL	0.09	0.15	0.18	0.07

Phase I Sampling Program - 11/10/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
t-1,2-Dichloroethylene	0.04	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.04	0.02
t-2-Butene	0.17	1.32	0.13	0.18	0.15	0.1	6	0.10	0.34	0.73	1.32	0.48
t-2-Hexene	0.04	<DL	<DL	0.02	<DL	<DL	6	<DL	0.01	0.02	0.04	0.02
t-2-Pentene	0.08	<DL	0.02	0.17	0.08	0.04	6	<DL	0.07	0.11	0.17	0.06
t-Butylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Tetrachloroethylene	0.04	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.04	0.02
Toluene	0.4	2.03	0.29	0.51	0.35	0.31	6	0.29	0.65	1.20	2.03	0.68
Trichloroethylene	0.08	0.04	<DL	<DL	<DL	0.04	6	<DL	0.03	0.05	0.08	0.03
Trichlorofluoromethane	0.47	0.33	0.35	0.36	0.37	0.33	6	0.33	0.37	0.41	0.47	0.05
Vinyl Acetate	<DL	<DL	0.37	<DL	0.6	<DL	6	<DL	0.16	0.37	0.60	0.26
Vinyl Bromide	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Vinyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,3-Dichloropropene (Total)	0.06	0.23	ND	ND	ND	ND	6	<DL	0.05	0.12	0.23	0.09
Xylene (Total)	0.33	0.21	0.12	0.3	0.18	0.11	6	0.11	0.21	0.28	0.33	0.09

<DL- Less than Detection Limit

Min- Minimum Value

Max- Maximum Value

NC- Not Calculated

STD- Standard Deviation

* Suspected Laboratory Contaminant

Phase I Sampling Program - 11/16/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
1,1,1-Trichloroethane	0.04	0.03	<DL	0.03	0.03	0.03	6	<DL	0.03	0.04	0.04	0.01
1,1,2,2-Tetrachloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,1,2-Trichloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,1-Dichloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,1-Dichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2,3-Trimethylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2,4-Trichlorobenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2,4-Trimethylbenzene	0.02	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	0.02	NC
1,2-Dibromoethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2-Dichloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,2-Dichloropropane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,3,5-Trimethylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,3-Butadiene	0.09	0.05	0.05	0.09	0.07	0.03	6	0.03	0.06	0.08	0.09	0.02
1,4-Dioxane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Butanol	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Decene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Heptene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Hexene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Methylcyclohexene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Nonene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Octene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Propanol	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1-Undecene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2,2,3-Trimethylpentane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2,2,4-Trimethylpentane	0.04	0.01	0.01	0.02	0.03	0.02	6	0.01	0.02	0.03	0.04	0.01
2,2,5-Trimethylhexane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2,3,4-Trimethylpentane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2,3-Dimethylbutane	0.03	0.01	<DL	0.02	0.02	0.02	6	<DL	0.02	0.03	0.03	0.01
2,3-Dimethylpentane	0.02	<DL	<DL	<DL	<DL	0.02	6	<DL	0.01	0.02	0.02	0.01
2,4,4-Trimethyl-1-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2,4-4-Trimethyl-2-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2,4-Dimethylpentane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2,5-Dimethylhexane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2-Ethyl-1-Butene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2-Methyl-1-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2-Methyl-2-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2-Methylheptane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
2-Propanol	<DL	<DL	<DL	2.29	0.8	1.37	6	<DL	0.74	1.51	2.29	0.94
3-Methyl-1-Butene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
3-Methylheptane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
3-Methylhexane	0.03	0.01	<DL	0.02	0.02	0.02	6	<DL	0.02	0.03	0.03	0.01
3-Methylpentane	0.04	0.01	0.01	0.03	0.03	0.02	6	0.01	0.02	0.03	0.04	0.01
4-Methyl-1-Pentene	0.03	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.03	0.01
4-Nonene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Acetaldehyde	1.47	1.27	0.99	1.88	1.13	1.78	6	0.99	1.42	1.71	1.88	0.36
Acetone (+)	2.67	2.05	1.03	0.83	0.55	2.12	6	0.55	1.54	2.24	2.67	0.85
Acetonitrile	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Acetylene	0.67	0.47	0.49	0.62	0.59	0.43	6	0.43	0.55	0.62	0.67	0.10
Acrylonitrile	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Allyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
a-Pinene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Benzaldehyde	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Benzene	0.19	0.09	0.11	0.15	0.15	0.15	6	0.09	0.14	0.17	0.19	0.04
Benzyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
b-Pinene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromochloromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromodichloromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromoform	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Bromomethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Butyl Acrylate	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Butyraldehyde	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
c-1,2-Dichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
c-2-Butene	0.03	<DL	<DL	0.05	0.07	0.17	6	<DL	0.05	0.11	0.17	0.06
c-2-Hexene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
c-2-Octene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC

Phase I Sampling Program - 11/16/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
c-2-Pentene	0.02	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	0.02	NC
c-3-Hexene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
c-3-Methyl-2-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
c-4-Methyl-2-Pentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Carbon Tetrachloride	0.12	0.12	0.12	0.13	0.13	0.13	6	0.12	0.13	0.13	0.13	0.01
Chlorobenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Chlorodifluoromethane	0.4	0.27	0.28	0.35	0.27	0.23	6	0.23	0.30	0.35	0.40	0.06
Chloroethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Chloroform	0.02	0.01	<DL	0.02	0.02	0.02	6	<DL	0.02	0.02	0.02	0.01
Chloromethane	0.53	0.5	0.5	0.52	0.47	0.49	6	0.47	0.50	0.52	0.53	0.02
Chloroprene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Cyclohexane	<DL	<DL	<DL	<DL	<DL	0.02	6	<DL	NC	NC	0.02	NC
Cyclohexene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Cyclopentane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Cyclopentene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Dibromochloromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Dichlorodifluoromethane	0.68	0.67	0.64	0.64	0.63	0.64	6	0.63	0.65	0.67	0.68	0.02
Dichlorofluoromethane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Diethyl Ether	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Epichlorohydrin	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Ethane	3.67	3.17	3.6	3.9	4.73	3.26	6	3.17	3.72	4.18	4.73	0.56
Ethanol	<DL	<DL	<DL	<DL	<DL	1.42	6	<DL	0.24	0.71	1.42	0.58
Ethylbenzene	0.03	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.03	0.01
Ethylene	1.18	0.93	1.18	1.23	1.18	0.95	6	0.93	1.11	1.22	1.23	0.13
Freon 113	0.1	0.1	0.1	0.09	0.09	0.09	6	0.09	0.10	0.10	0.10	0.01
Freon 114	0.01	0.01	<DL	0.01	<DL	0.01	6	<DL	0.01	0.01	0.01	0.01
Halocarbon 134A	1.36	0.48	0.09	0.28	0.15	1.18	6	0.09	0.59	1.04	1.36	0.55
Heptanal	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Hexachloro-1,3-Butadiene*	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Hexanal	0.05	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.03	0.05	0.02
Indan	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Indene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Isobutane	0.75	0.23	0.25	0.32	0.32	0.3	6	0.23	0.36	0.52	0.75	0.19
Isobutene + 1-Butene	0.19	0.03	<DL	0.1	0.09	0.16	6	<DL	0.10	0.15	0.19	0.07
Isobutylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Isoheptane	0.03	0.01	<DL	0.02	0.03	0.02	6	<DL	0.02	0.03	0.03	0.01
Isohexane	0.12	0.04	0.05	0.07	0.08	0.05	6	0.04	0.07	0.09	0.12	0.03
Isoprene	0.11	<DL	<DL	<DL	<DL	0.03	6	<DL	0.02	0.06	0.11	0.04
m-Dichlorobenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
m-Diethylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Methyl ethyl ketone	<DL	<DL	<DL	<DL	0.98	<DL	6	<DL	0.16	0.49	0.98	0.40
Methyl t-Butylether	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Methylcyclohexane	0.06	<DL	<DL	<DL	<DL	0.04	6	<DL	0.02	0.04	0.06	0.03
Methylcyclopentane	0.02	<DL	<DL	0.02	0.02	0.02	6	<DL	0.01	0.02	0.02	0.01
Methylene Chloride	0.47	0.24	0.13	0.51	0.13	0.41	6	0.13	0.32	0.45	0.51	0.17
Methylisobutylketone	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Naphthalene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
n-Butane	1.87	0.71	0.64	0.86	1.07	1.39	6	0.64	1.09	1.47	1.87	0.47
n-Butylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
n-Decane	0	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Neohexane	0.03	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.03	0.01
Neopentane	0.01	<DL	<DL	<DL	<DL	0	6	<DL	NC	NC	0.01	NC
n-Heptane	0.04	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.04	0.02
n-Hexane	0.07	0.02	0.03	0.05	0.07	0.04	6	0.02	0.05	0.06	0.07	0.02
n-Nonane	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
n-Octane	0.02	<DL	<DL	<DL	<DL	0.01	6	<DL	0.01	0.01	0.02	0.01
n-Pentane	0.37	0.17	0.18	0.37	0.22	0.18	6	0.17	0.25	0.33	0.37	0.10
n-Propylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
n-Undecane	0.01	<DL	<DL	<DL	<DL	0.01	6	<DL	NC	NC	0.01	NC
o-Dichlorobenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
p-Dichlorobenzene	0.78	0.16	0.08	0.11	0.06	0.46	6	0.06	0.28	0.51	0.78	0.29
p-Diethylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
p-Isopropyltoluene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Propane	2.52	2.43	1.38	1.76	2.27	2.75	6	1.38	2.19	2.61	2.75	0.52
Propylene	0.56	0.33	0.4	0.38	0.94	0.14	6	0.14	0.46	0.68	0.94	0.27
Styrene	0.03	0.03	<DL	0.09	<DL	<DL	6	<DL	0.03	0.05	0.09	0.04

Phase I Sampling Program - 11/16/02 Sampling Program

	Norco Phase I Community Sampling Project Individual Site Sampling Results (ppb)						Norco Phase I Community Sampling Project Individual Site Sampling Statistics (ppb)					
Compound Name	**Site 1**	**Site 2**	**Site 3**	**Site 4**	**Site 5**	**Site 6**	**Sample Count**	**Min (ppb)**	**Average (ppb)**	**Upper 95% Confidence Limit, UCL (ppb)**	**Max (ppb)**	**STD (ppb)**
t-1,2-Dichloroethylene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
t-2-Butene	0.04	<DL	<DL	0.11	0.14	0.26	6	<DL	0.09	0.17	0.26	0.10
t-2-Hexene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
t-2-Pentene	0.03	<DL	<DL	<DL	<DL	<DL	6	<DL	0.01	0.02	0.03	0.01
t-Butylbenzene	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Tetrachloroethylene	0.49	0.14	0.05	0.08	0.03	0.42	6	0.03	0.20	0.37	0.49	0.20
Toluene	0.69	0.81	0.2	0.54	0.49	2.17	6	0.20	0.82	1.38	2.17	0.69
Trichloroethylene	0.7	0.22	0.08	0.13	0.07	0.71	6	0.07	0.32	0.57	0.71	0.30
Trichlorofluoromethane	0.35	0.34	0.34	0.33	0.32	0.34	6	0.32	0.34	0.35	0.35	0.01
Vinyl Acetate	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Vinyl Bromide	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
Vinyl Chloride	<DL	<DL	<DL	<DL	<DL	<DL	6	<DL	NC	NC	<DL	NC
1,3-Dichloropropene (Total)	ND	ND	ND	ND	ND	ND	6	<DL	NC	NC	<DL	NC
Xylene (Total)	0.09	ND	ND	0.03	0.03	ND	6	<DL	0.03	0.05	0.09	0.04

<DL- Less than Detection Limit

Min- Minimum Value

Max- Maximum Value

NC- Not Calculated

STD- Standard Deviation

* Suspected Laboratory Contaminant