LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN100720\
Data File : VN063951.D
Acq On : 7 Oct 2020 18:31
Operator : JC/MD
Sample : L4250-05
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-10

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.997	70	80	90	rBV	54558	78249	4.70%	0.399%
2	2.161	121	131	143	rVB	66939	101620	6.11%	0.519%
3	2.242	144	156	165	rBV	33404	56124	3.37%	0.286%
4	2.428	199	214	225	rBV9	18582	41553	2.50%	0.212%
5	2.775	308	322	338	rBV	131791	272766	16.39%	1.392%
6	3.406	507	518	537	rVB4	7505	19381	1.16%	0.099%
7	3.531	541	557	572	rBV3	14106	35679	2.14%	0.182%
8	3.762	605	629	654	rBV3	105206	331922	19.95%	1.694%
9	4.016	690	708	724	rBV2	13790	40207	2.42%	0.205%
10	4.486	829	854	874	rBV3	225166	792702	47.63%	4.045%
11	4.592	876	887	912	rVB4	52151	171102	10.28%	0.873%
12	5.010	998	1017	1037	rVB5	14963	51501	3.09%	0.263%
13	6.016	1315	1330	1348	rBV9	6168	17947	1.08%	0.092%
14	6.315	1406	1423	1435	rBV7	7851	25877	1.55%	0.132%
15	6.476	1454	1473	1491	rBV3	33149	101841	6.12%	0.520%
16	6.589	1491	1508	1519	rVV6	18496	54118	3.25%	0.276%
17	6.669	1520	1533	1549	rVB8	15781	45448	2.73%	0.232%
18	7.351	1731	1745	1761	rVB5	16490	46796	2.81%	0.239%
19	7.550	1786	1807	1818	rBV	209692	523857	31.48%	2.673%
20	7.627	1819	1831	1844	rVV	368277	905060	54.39%	4.618%
21	7.704	1844	1855	1876	rVB3	112948	302805	18.20%	1.545%
22	7.897	1899	1915	1928	rBV5	20241	50715	3.05%	0.259%
23	7.991	1928	1944	1966	rVV2	266565	658588	39.57%	3.360%
24	8.110	1966	1981	1989	rVV5	20890	58137	3.49%	0.297%
25	8.174	1990	2001	2020	rVV3	31917	99167	5.96%	0.506%
26	8.264	2020	2029	2043	rVB10	10993	23359	1.40%	0.119%
27	8.550	2101	2118	2142	rBV	557981	1187749	71.37%	6.061%
28	8.788	2181	2192	2199	rBV6	12156	24072	1.45%	0.123%
29	9.013	2248	2262	2272	rBV6	26777	64018	3.85%	0.327%
30	9.328	2341	2360	2373	rBV5	14369	35233	2.12%	0.180%
31	9.489	2399	2410	2425	rVB3	31053	65162	3.92%	0.332%
32	9.624	2429	2452	2465	rBV4	61898	170748	10.26%	0.871%
33	9.814	2498	2511	2526	rBV9	23507	54656	3.28%	0.279%
34	10.061	2573	2588	2603	rBV	912431	1664154	100.00%	8.491%

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Operator : JC/MD
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Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-10

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
Title : SW846 8260

35	10.273	2643	2654	2666	rVB	139001	250299	15.04%	1.277%
36	10.351	2666	2678	2688	rVV2	144125	251518	15.11%	1.283%
37	10.402	2688	2694	2705	rVB5	14083	26991	1.62%	0.138%
38	10.492	2705	2722	2738	rBV3	55916	122107	7.34%	0.623%
39	10.572	2738	2747	2759	rVB6	12069	25278	1.52%	0.129%
40	10.765	2790	2807	2818	rBV2	45792	85863	5.16%	0.438%
41	10.891	2837	2846	2856	rVB6	18485	32598	1.96%	0.166%
42	11.090	2899	2908	2921	rVB6	12065	24523	1.47%	0.125%
43	11.228	2942	2951	2961	rVB	31239	50211	3.02%	0.256%
44	11.379	2984	2998	3016	rBV	917396	1601139	96.21%	8.170%
45	11.486	3020	3031	3048	rVV	88433	183098	11.00%	0.934%
46	11.598	3052	3066	3086	rVB	147499	300955	18.08%	1.536%
47	11.923	3157	3167	3179	rVB	110211	182834	10.99%	0.933%
48	12.096	3211	3221	3231	rBV	12294	24267	1.46%	0.124%
49	12.164	3236	3242	3250	rVV	10963	17553	1.05%	0.090%
50	12.222	3250	3260	3270	rVV2	25294	43665	2.62%	0.223%
51	12.376	3295	3308	3325	rBV	744682	1281135	76.98%	6.537%
52	12.566	3362	3367	3378	rVB3	18090	27992	1.68%	0.143%
53	12.633	3380	3388	3392	rBV	39063	59295	3.56%	0.303%
54	12.707	3405	3411	3424	rVB3	22849	39033	2.35%	0.199%
55	12.868	3452	3461	3470	rBV2	48191	78182	4.70%	0.399%
56	13.016	3498	3507	3519	rVB	151138	248749	14.95%	1.269%
57	13.135	3532	3544	3556	rVB4	44268	100188	6.02%	0.511%
58	13.212	3561	3568	3577	rVB9	13400	21579	1.30%	0.110%
59	13.318	3588	3601	3623	rBV	870474	1549599	93.12%	7.907%
60	13.418	3624	3632	3641	rVB	37743	60583	3.64%	0.309%
61	13.489	3643	3654	3660	rBV	293213	489191	29.40%	2.496%
62	13.566	3671	3678	3693	rVB2	113463	186197	11.19%	0.950%
63	13.649	3694	3704	3712	rVV2	65360	101556	6.10%	0.518%
64	13.704	3712	3721	3734	rVB5	43735	94297	5.67%	0.481%
65	13.785	3734	3746	3761	rBV4	100961	214425	12.88%	1.094%
66	13.865	3761	3771	3782	rVB	55067	86683	5.21%	0.442%
67	13.971	3791	3804	3816	rBV	797754	1318715	79.24%	6.729%
68	14.106	3835	3846	3854	rBV	69342	107063	6.43%	0.546%
69	14.186	3856	3871	3877	rVV4	64582	140548	8.45%	0.717%
70	14.222	3877	3882	3890	rVV	67209	105350	6.33%	0.538%
71	14.267	3890	3896	3901	rVV3	51593	84005	5.05%	0.429%

LSC Area Percent Report

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Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-10

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
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Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
Title : SW846 8260

72	14.299	3902	3906	3916	rVB2	62422	86586	5.20%	0.442%
73	14.402	3929	3938	3947	rBV2	43223	70512	4.24%	0.360%
74	14.460	3947	3956	3965	rBV4	69821	121600	7.31%	0.620%
75	14.566	3977	3989	4001	rBV8	29166	67169	4.04%	0.343%
76	14.633	4001	4010	4019	rBV4	102211	176764	10.62%	0.902%
77	14.714	4029	4035	4042	rVB2	26412	36809	2.21%	0.188%
78	14.826	4060	4070	4083	rBV4	83584	171590	10.31%	0.876%
79	14.916	4085	4098	4113	rVB2	208869	418179	25.13%	2.134%
80	15.103	4141	4156	4167	rVB	96302	192713	11.58%	0.983%
81	15.161	4168	4174	4183	rBV4	17038	24530	1.47%	0.125%
82	15.569	4295	4301	4321	rVB4	18025	45866	2.76%	0.234%
83	15.665	4321	4331	4340	rBV2	46086	90840	5.46%	0.464%
84	15.739	4344	4354	4363	rVB5	27003	61978	3.72%	0.316%
85	15.868	4383	4394	4403	rBV5	11222	23147	1.39%	0.118%
86	16.061	4445	4454	4464	rVB2	31064	60351	3.63%	0.308%
87	16.122	4464	4473	4487	rVB5	36955	73996	4.45%	0.378%
88	16.315	4520	4533	4551	rBV7	32899	85846	5.16%	0.438%

Sum of corrected areas: 19598053

LSC Report - Integrated Chromatogram

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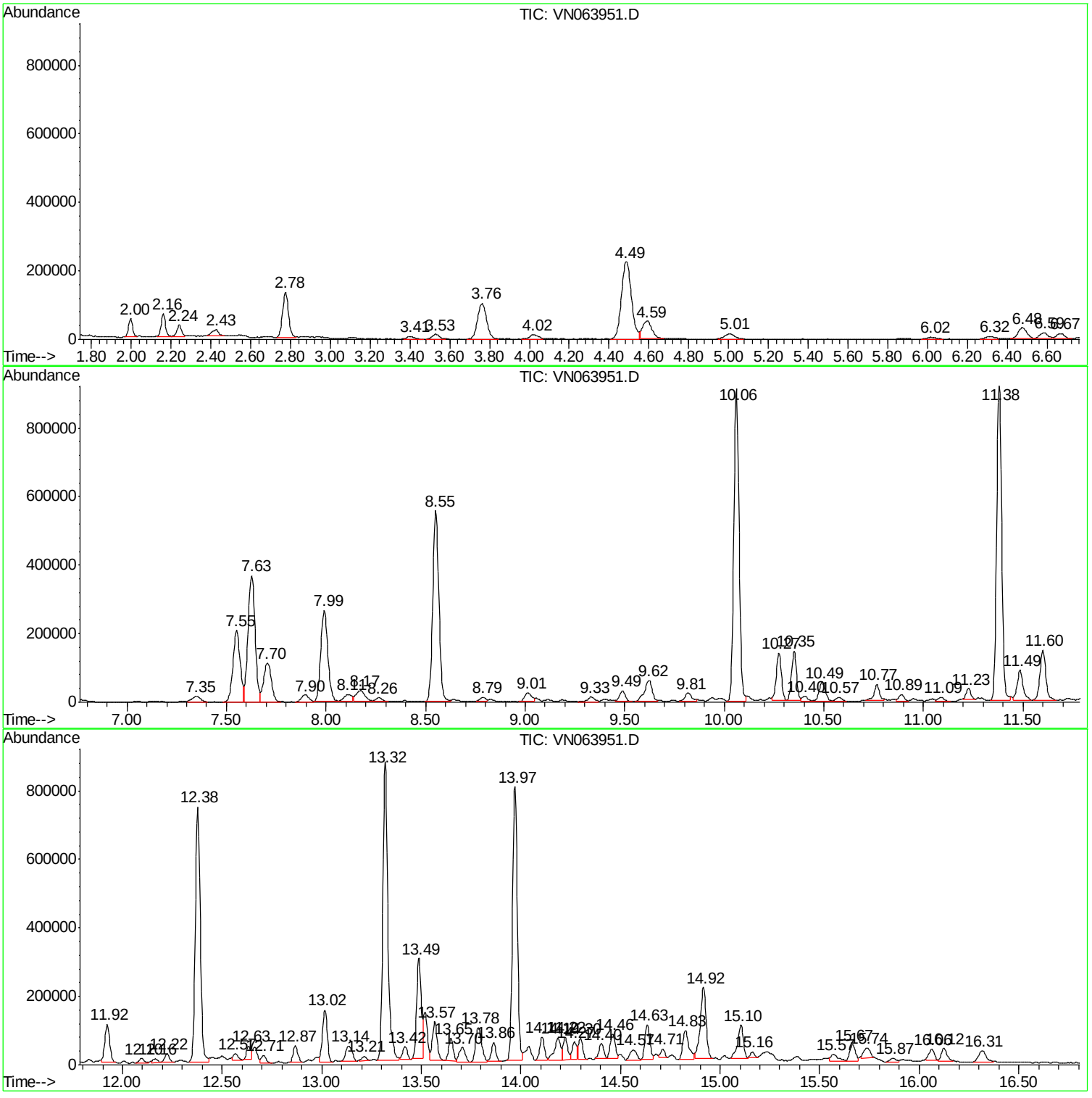
Sample : L4250-05
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100720\
Data File : VN063951.D
Acq On : 7 Oct 2020 18:31
Operator : JC/MD
Sample : L4250-05
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ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
Quant Title : SW846 8260

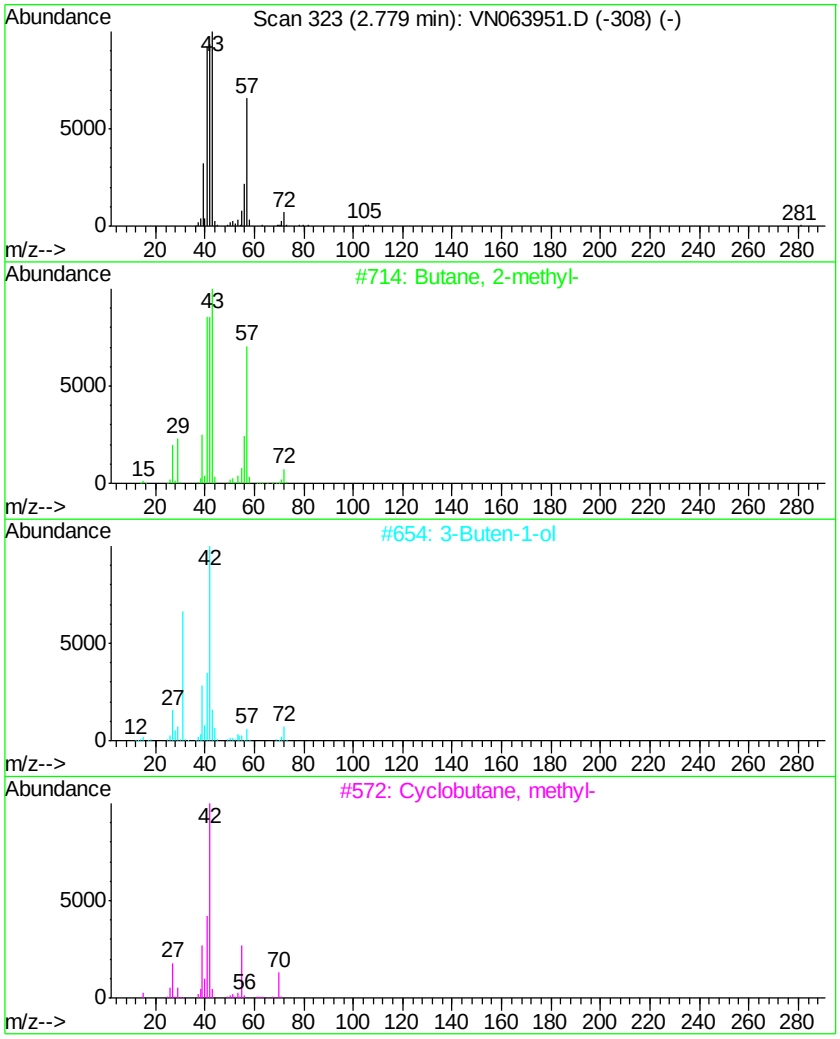
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

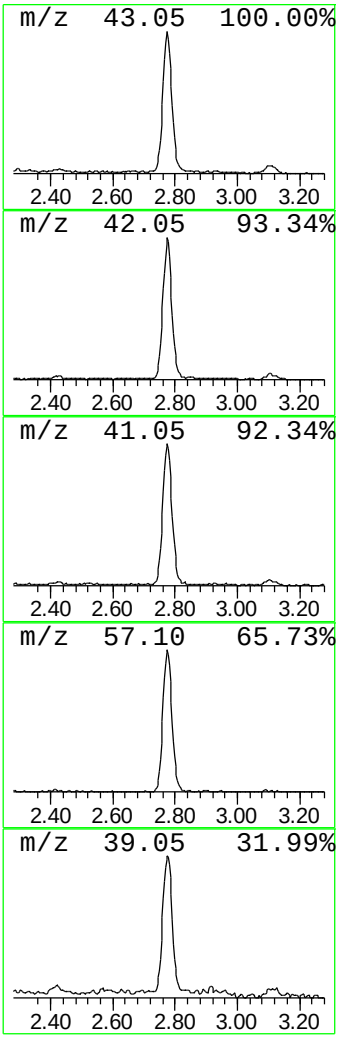
****

Peak Number 1 Butane, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.78	15.07 ug/l	272766	Pentafluorobenzene	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	91
2	3-Buten-1-ol	72	C4H8O	000627-27-0	47
3	Cyclobutane, methyl-	70	C5H10	000598-61-8	38
4	Pentane	72	C5H12	000109-66-0	36
5	2(3H)-Furanone, dihydro-4-methyl-	100	C5H8O2	001679-49-8	33





Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100720\
Data File : VN063951.D
Acq On : 7 Oct 2020 18:31
Operator : JC/MD
Sample : L4250-05
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
Quant Title : SW846 8260

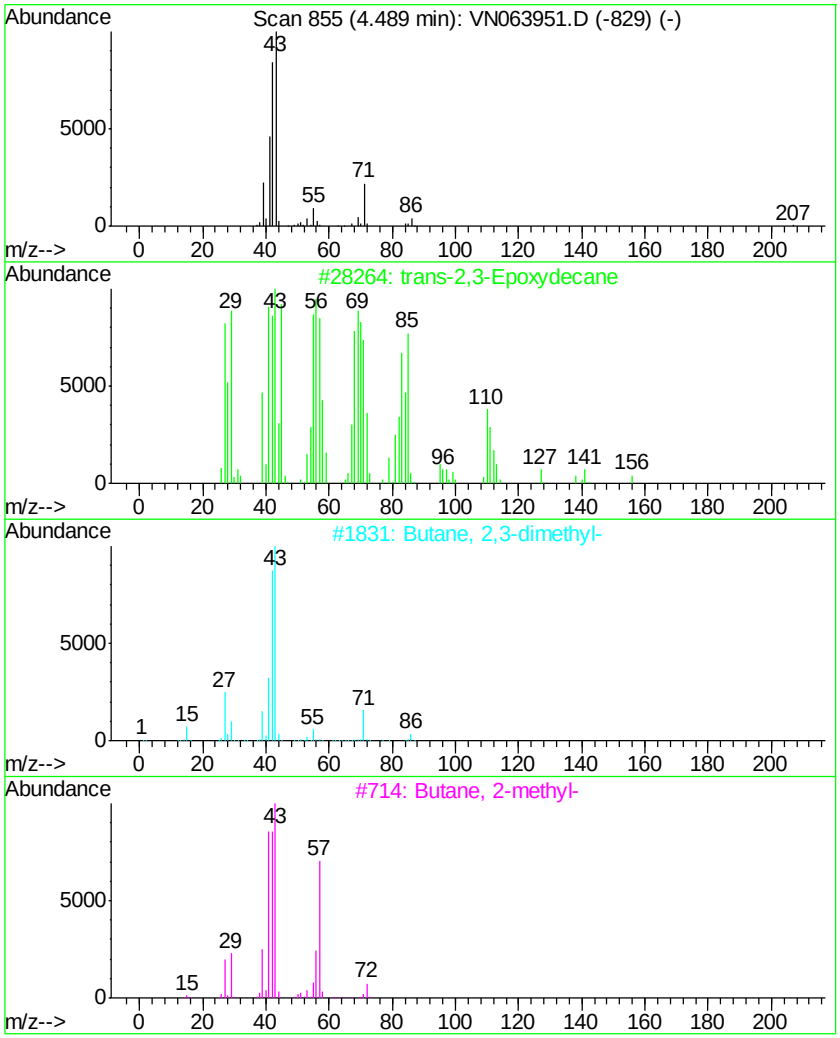
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

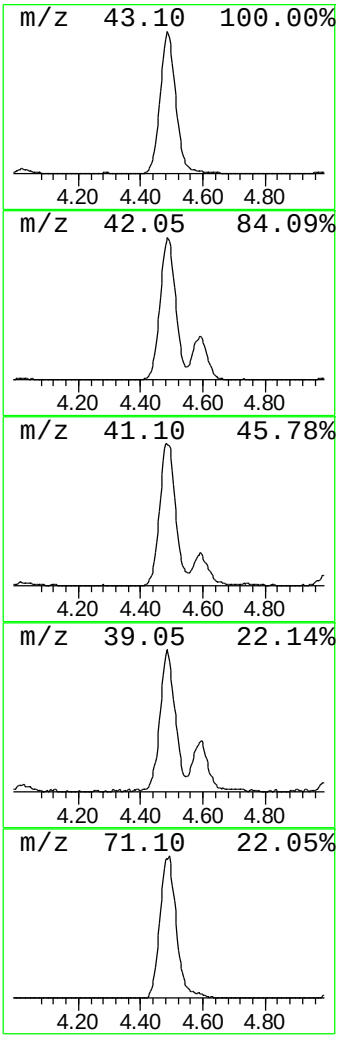
****

Peak Number 2 trans-2,3-Epoxydecane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.49	43.79 ug/l	792702	Pentafluorobenzene	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	trans-2,3-Epoxydecane	156	C10H20O	054125-39-2	87
2	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	86
3	Butane, 2-methyl-	72	C5H12	000078-78-4	38
4	1-Tridecene	182	C13H26	002437-56-1	35
5	Pentane, 2-methyl-	86	C6H14	000107-83-5	23





Library Search Compound Report

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ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
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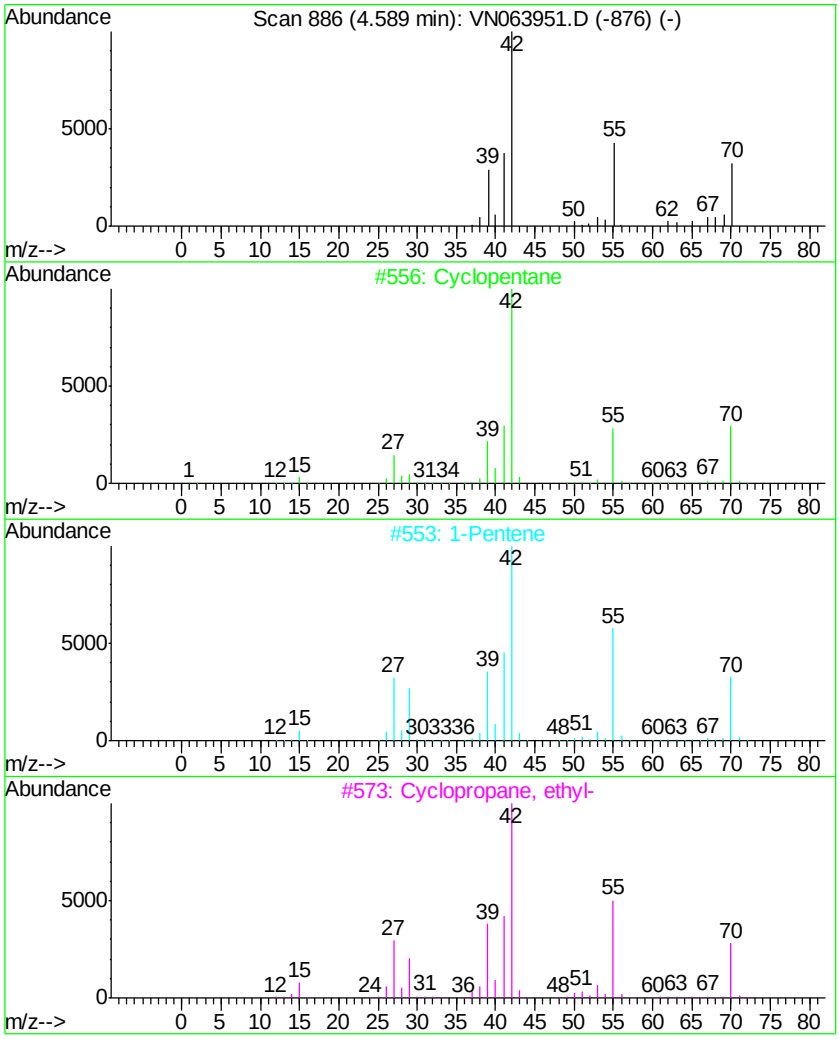
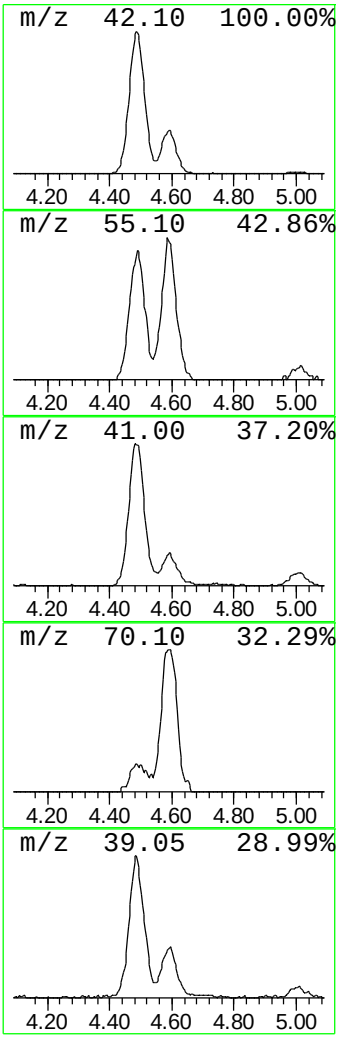
****

Peak Number 3 Cyclopentane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.59	9.45 ug/l	171102	Pentafluorobenzene	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane	70	C5H10	000287-92-3	64
2	1-Pentene	70	C5H10	000109-67-1	64
3	Cyclopropane, ethyl-	70	C5H10	001191-96-4	9
4	2(3H)-Furanone, dihydro-3,5-dime...	114	C6H10O2	005145-01-7	9
5	Cyclobutane, methyl-	70	C5H10	000598-61-8	9





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ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
Quant Title : SW846 8260

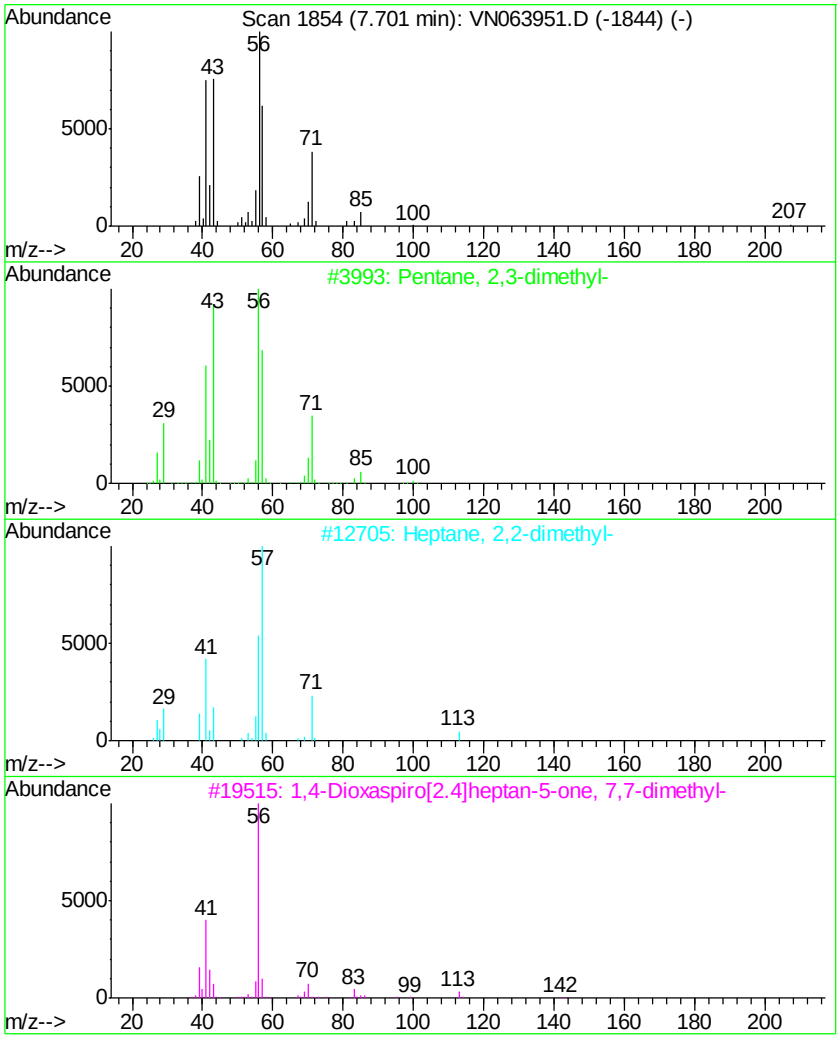
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

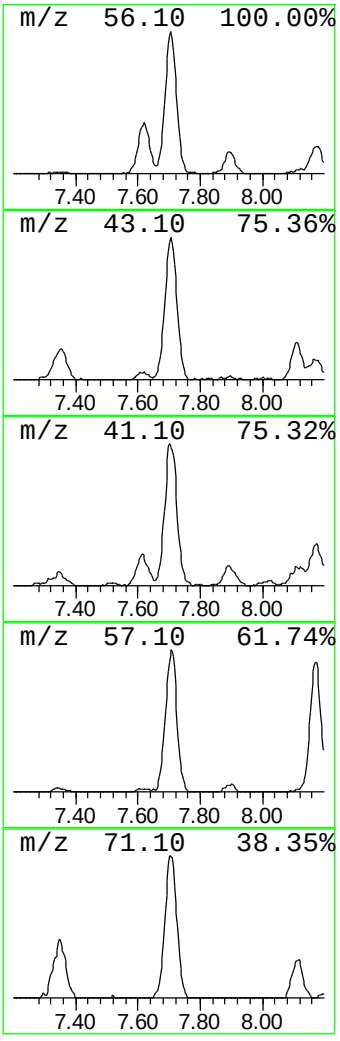
****

Peak Number 4 Pentane, 2,3-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.70	16.73 ug/l	302805	Pentafluorobenzene	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	90
2	Heptane, 2,2-dimethyl-	128	C9H20	001071-26-7	47
3	1,4-Dioxaspiro[2.4]heptan-5-one,...	142	C7H10O3	126091-79-0	40
4	1-Hexene, 2-methyl-	98	C7H14	006094-02-6	38
5	1-Hexene, 2,5-dimethyl-	112	C8H16	006975-92-4	38





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Sample : L4250-05
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
Quant Title : SW846 8260

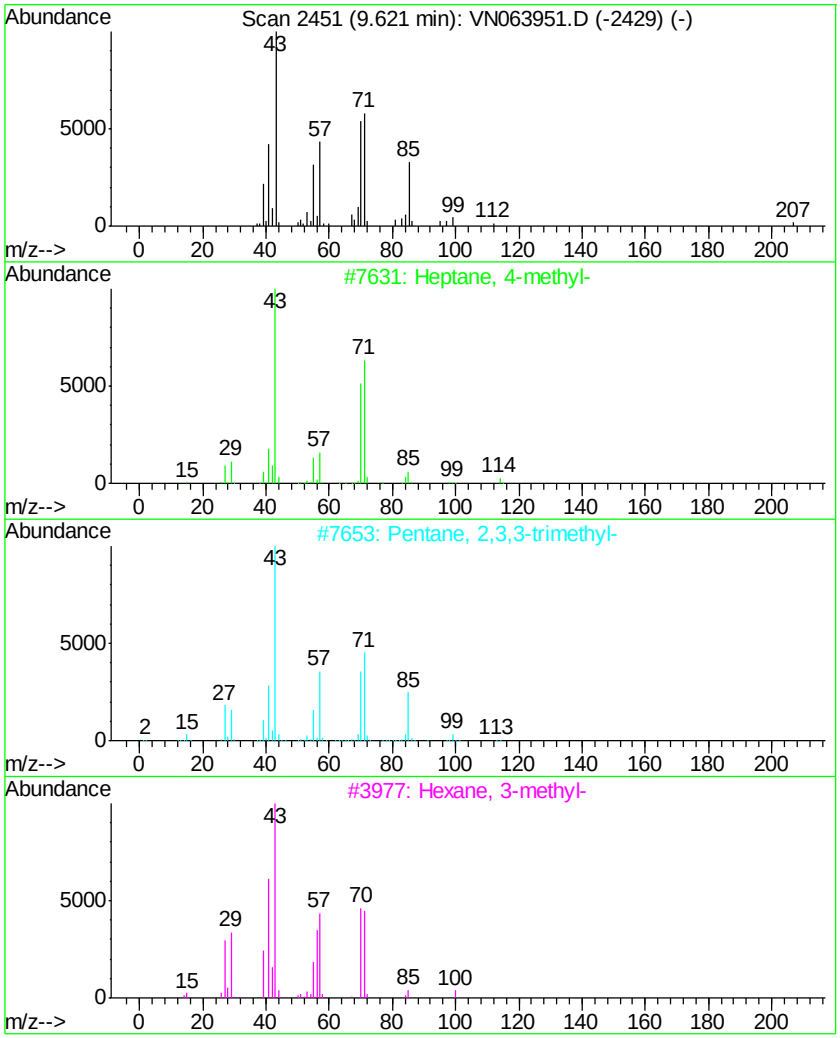
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

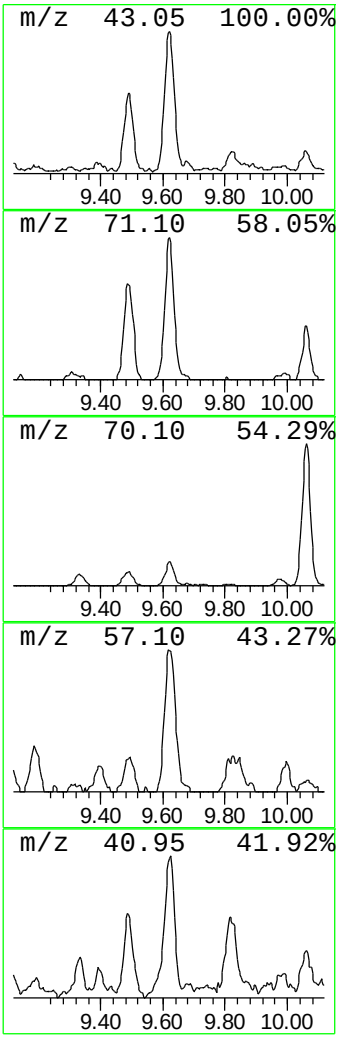
****

Peak Number 5 Heptane, 4-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.62	7.19 ug/l	170748	1,4-Difluorobenzene	8.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 4-methyl-	114	C8H18	000589-53-7	64
2	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	64
3	Hexane, 3-methyl-	100	C7H16	000589-34-4	59
4	1-Heptene, 4-methyl-	112	C8H16	013151-05-8	55
5	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	53





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Operator : JC/MD
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Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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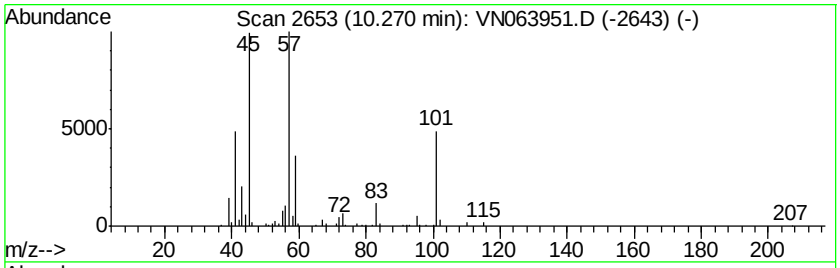
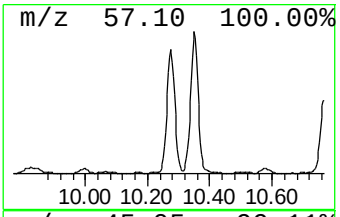
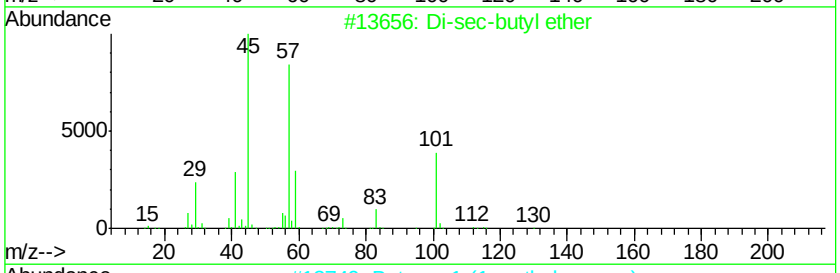
Peak Number 6 Di-sec-butyl ether Concentration Rank 9

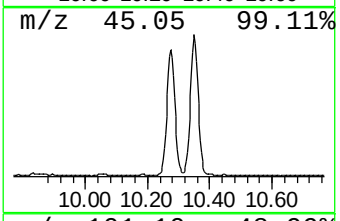
R.T.	EstConc	Area	Relative to ISTD	R.T.
10.27	7.82 ug/l	250299	Chlorobenzene-d5	11.38

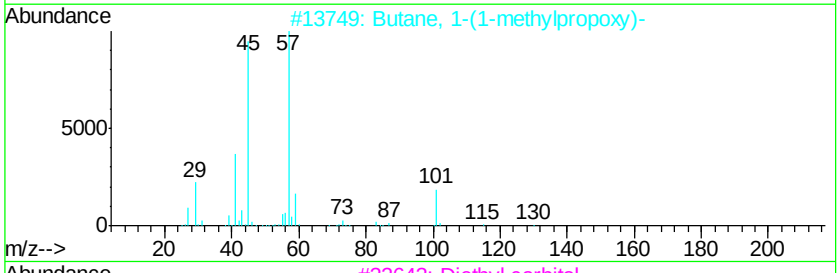
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Di-sec-butyl ether	130	C8H18O	006863-58-7	90
2	Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	56
3	Diethyl carbitol	162	C8H18O3	000112-36-7	43
4	3-Pentanol, 3-ethyl-2-methyl-	130	C8H18O	000597-05-7	36
5	Boronic acid, ethyl-, diethyl ester	130	C6H15BO2	053907-92-9	36

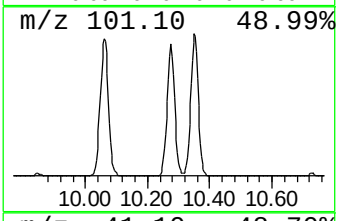


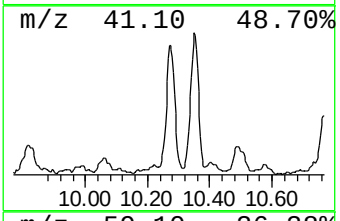


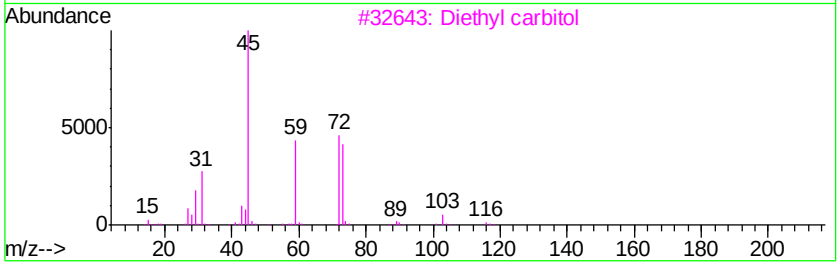


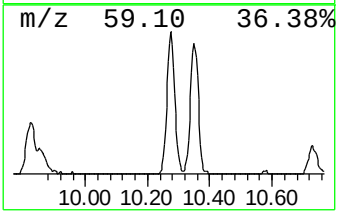












Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN100720\
Data File : VN063951.D
Acq On : 7 Oct 2020 18:31
Operator : JC/MD
Sample : L4250-05
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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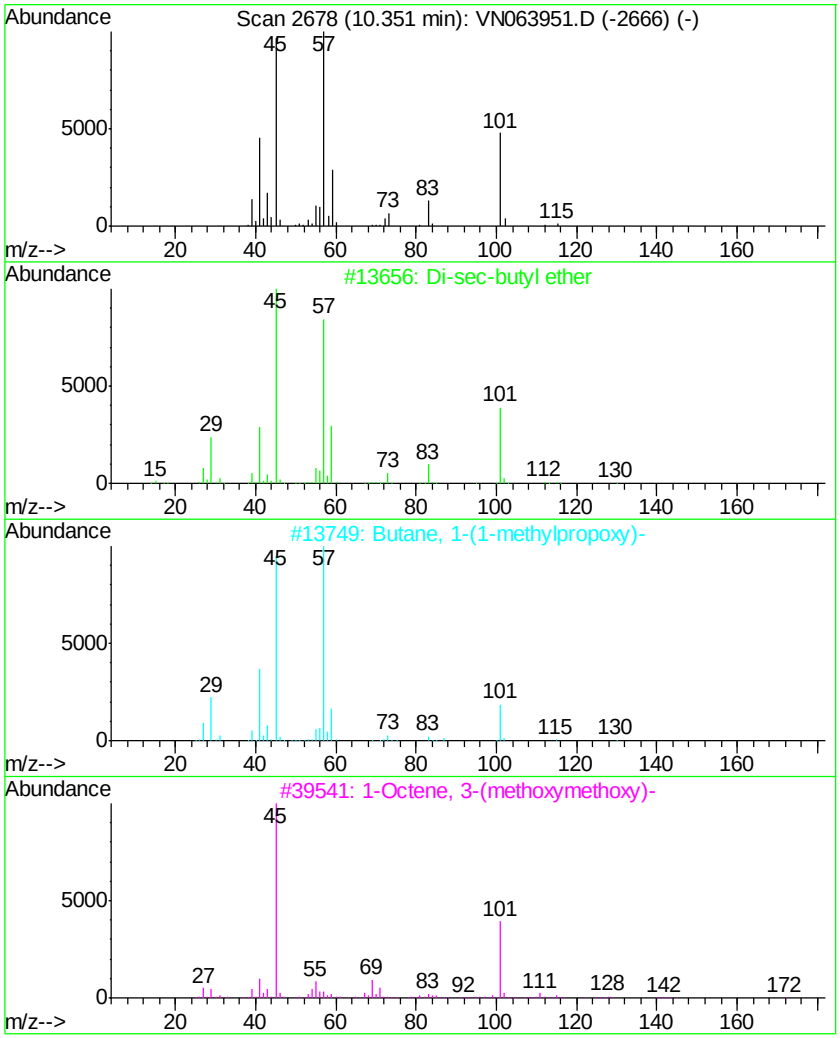
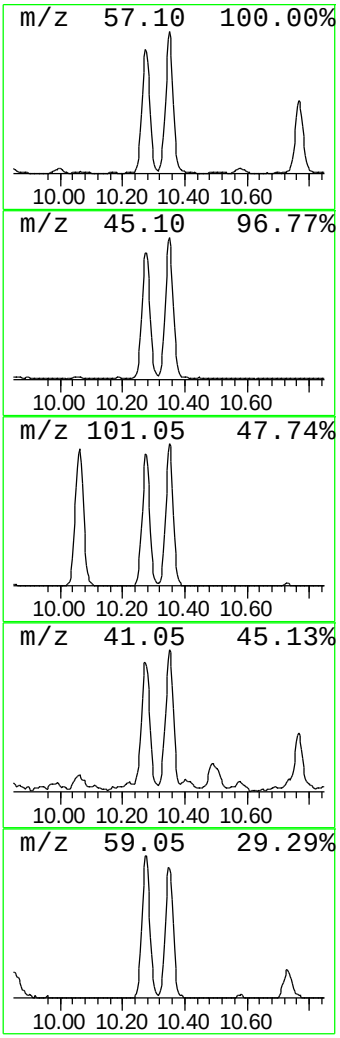
Peak Number 7 Butane, 1-(1-methylpropoxy)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.35	7.85 ug/l	251518	Chlorobenzene-d5	11.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Di-sec-butyl ether	130	C8H18O	006863-58-7	90
2	Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	53
3	1-Octene, 3-(methoxymethoxy)-	172	C10H20O2	088738-34-5	40
4	3-Pentanol, 3-ethyl-2-methyl-	130	C8H18O	000597-05-7	36
5	Fructofuranose, 2,6-anhydro-1,3,...	204	C9H16O5	004451-11-0	33





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN100720\
Data File : VN063951.D
Acq On : 7 Oct 2020 18:31
Operator : JC/MD
Sample : L4250-05
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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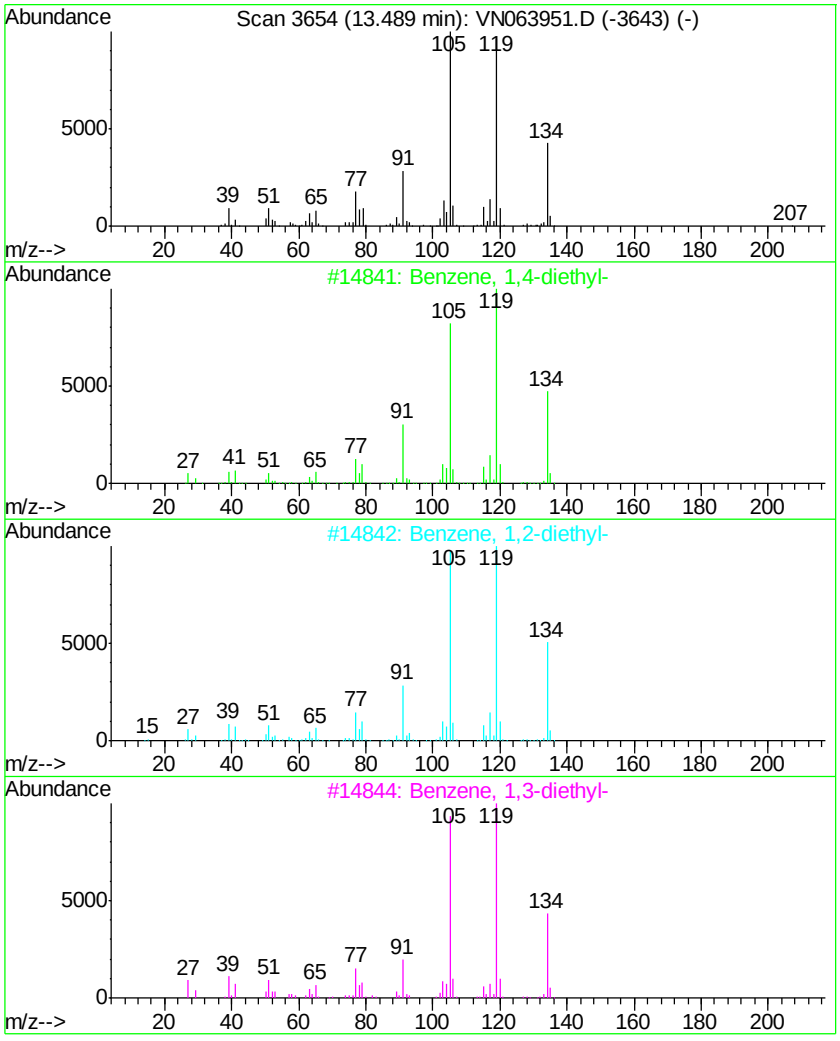
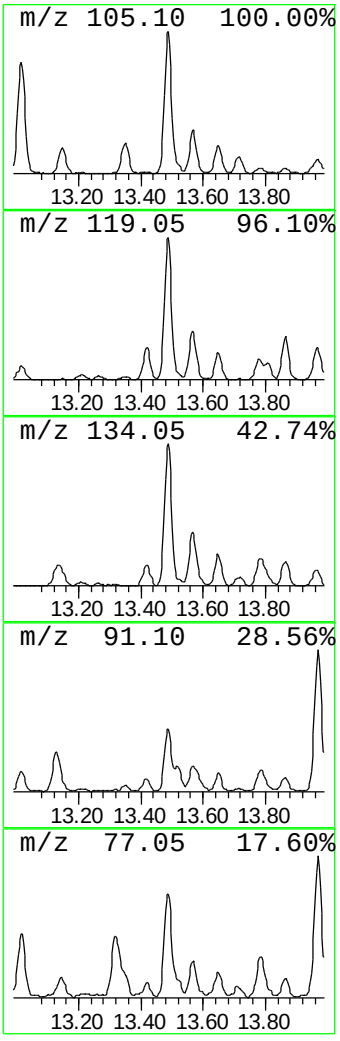
Peak Number 8 Benzene, 1,4-diethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.49	15.78 ug/l	489191	1,4-Dichlorobenzene-d4	13.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	97
2	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	96
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN100720\
Data File : VN063951.D
Acq On : 7 Oct 2020 18:31
Operator : JC/MD
Sample : L4250-05
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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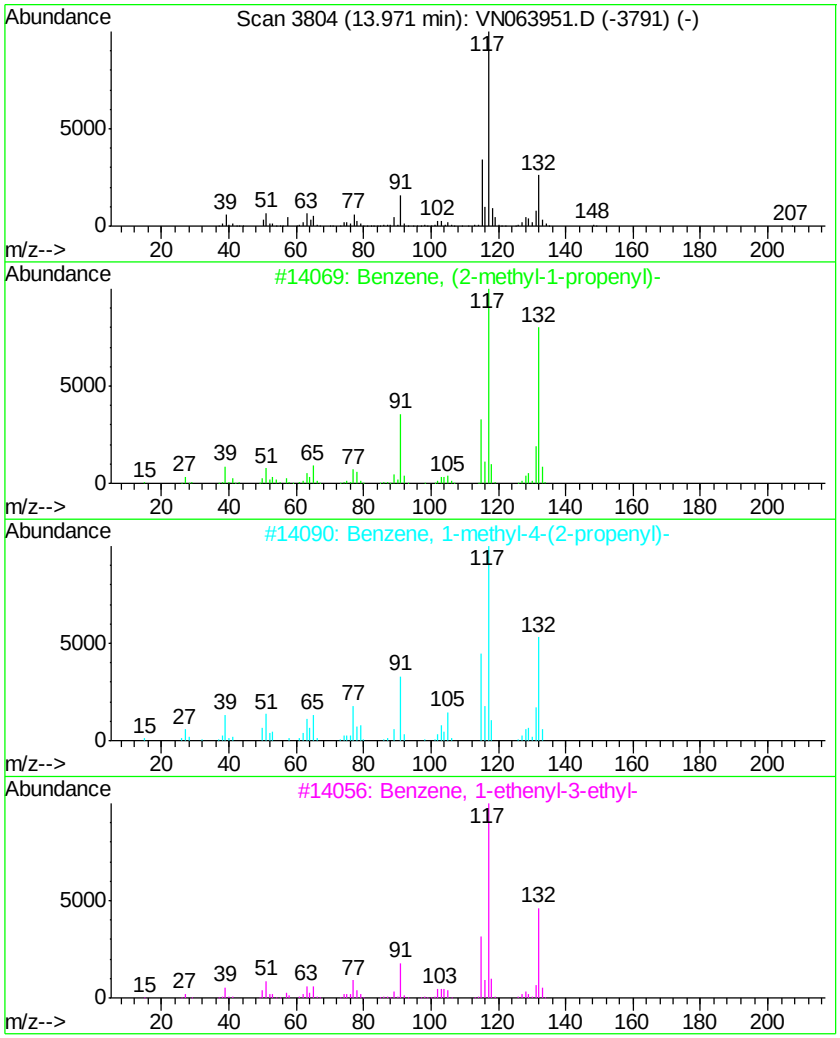
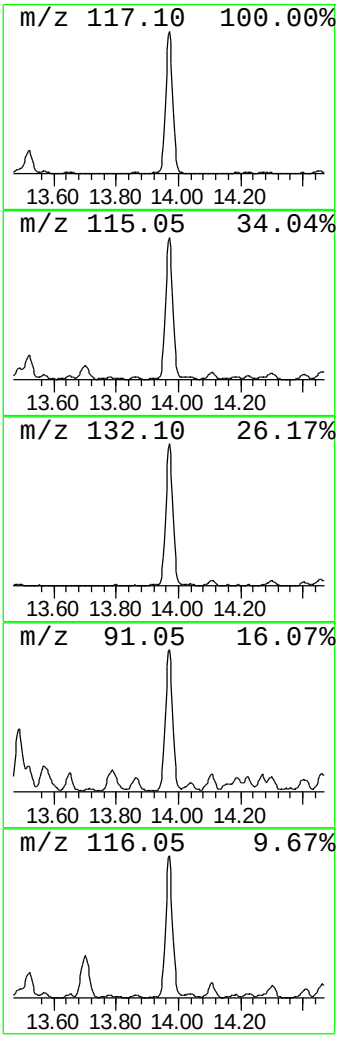
Peak Number 9 Benzene, (2-methyl-1-propen... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	42.55 ug/l	1318720	1,4-Dichlorobenzene-d4	13.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90
2	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	87
3	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87
5	Indan, 1-methyl-	132	C10H12	000767-58-8	87





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN100720\
Data File : VN063951.D
Acq On : 7 Oct 2020 18:31
Operator : JC/MD
Sample : L4250-05
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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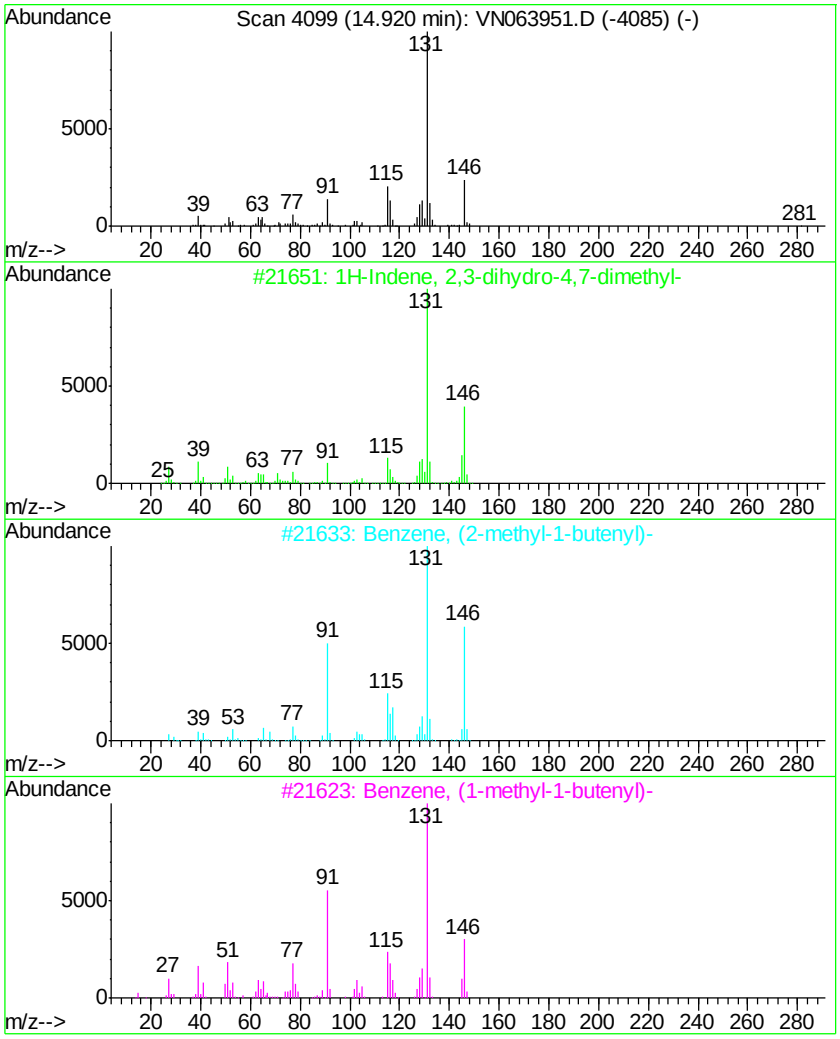
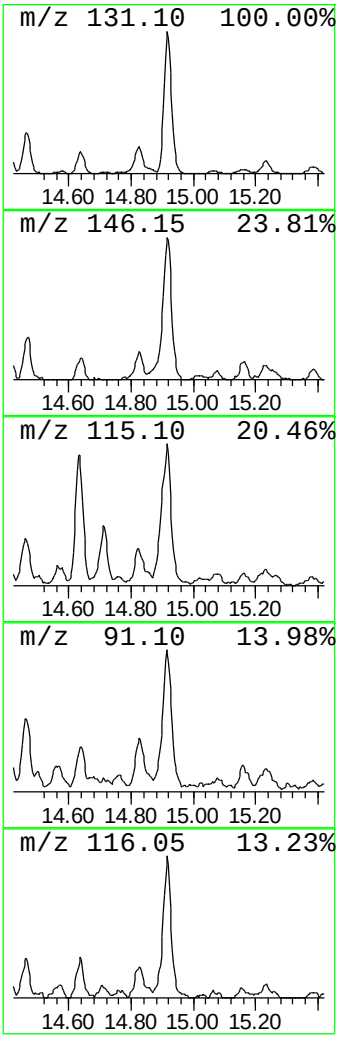
Peak Number 10 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.92	13.49 ug/l	418179	1,4-Dichlorobenzene-d4	13.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91
2	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	91
3	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	91
4	1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	90
5	1,3-Cyclopentadiene, 5-(trans-2-...	146	C11H14	079209-36-2	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN100720\
Data File : VN063951.D
Acq On : 7 Oct 2020 18:31
Operator : JC/MD
Sample : L4250-05
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-methyl-	2.78	15.1	ug/l	272766	1	7.63	905060	50.0
trans-2,3-Epoxyde...	4.49	43.8	ug/l	792702	1	7.63	905060	50.0
Cyclopentane	4.59	9.4	ug/l	171102	1	7.63	905060	50.0
Pentane, 2,3-dime...	7.70	16.7	ug/l	302805	1	7.63	905060	50.0
Heptane, 4-methyl-	9.62	7.2	ug/l	170748	2	8.55	1187750	50.0
Di-sec-butyl ether	10.27	7.8	ug/l	250299	3	11.38	1601140	50.0
Butane, 1-(1-meth...	10.35	7.8	ug/l	251518	3	11.38	1601140	50.0
Benzene, 1,4-diet...	13.49	15.8	ug/l	489191	4	13.32	1549600	50.0
Benzene, (2-methy...	13.97	42.5	ug/l	1318720	4	13.32	1549600	50.0
1H-Indene, 2,3-di...	14.92	13.5	ug/l	418179	4	13.32	1549600	50.0