LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN112420\
Data File : VN064807.D
Acq On : 24 Nov 2020 13:50
Operator : JC/MD
Sample : L4829-12
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-2

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\82N112320W.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	2.782	309	324	351	rBV	329615	670592	23.81%	1.575%
2	3.106	413	425	444	rVB2	63435	139498	4.95%	0.328%
3	3.524	538	555	573	rVB2	19790	49121	1.74%	0.115%
4	3.756	606	627	652	rBV5	46920	151428	5.38%	0.356%
5	4.486	831	854	858	rBV2	133772	354322	12.58%	0.832%
6	5.000	985	1014	1052	rBV2	333526	1193669	42.37%	2.803%
7	5.518	1152	1175	1199	rBV4	35983	112447	3.99%	0.264%
8	5.878	1267	1287	1310	rVB2	50188	152816	5.42%	0.359%
9	6.020	1311	1331	1350	rBV3	37004	119353	4.24%	0.280%
10	6.319	1402	1424	1446	rBV9	41066	146609	5.20%	0.344%
11	6.476	1456	1473	1486	rBV3	36692	105830	3.76%	0.249%
12	6.579	1487	1505	1525	rBV2	242128	738021	26.20%	1.733%
13	6.785	1552	1569	1586	rVB8	13386	44429	1.58%	0.104%
14	7.161	1663	1686	1704	rVB7	9052	35511	1.26%	0.083%
15	7.283	1706	1724	1732	rBV4	26846	74393	2.64%	0.175%
16	7.344	1733	1743	1762	rVB3	58486	158814	5.64%	0.373%
17	7.550	1785	1807	1814	rBV	123510	332806	11.81%	0.782%
18	7.621	1816	1829	1843	rVV6	325799	973713	34.57%	2.287%
19	7.708	1844	1856	1878	rVB3	177217	496374	17.62%	1.166%
20	7.836	1878	1896	1909	rBV2	170165	418179	14.85%	0.982%
21	8.000	1928	1947	1969	rBV3	693214	1757708	62.40%	4.128%
22	8.122	1969	1985	1994	rBV6	188232	505059	17.93%	1.186%
23	8.264	2022	2029	2050	rVB3	86054	207769	7.38%	0.488%
24	8.373	2050	2063	2091	rVB5	37183	118443	4.20%	0.278%
25	8.547	2098	2117	2139	rBV3	529216	1329793	47.21%	3.123%
26	8.643	2140	2147	2158	rVB4	35593	66012	2.34%	0.155%
27	8.711	2158	2168	2179	rVB3	48927	93928	3.33%	0.221%
28	8.788	2179	2192	2199	rBV2	70657	151596	5.38%	0.356%
29	9.016	2243	2263	2264	rBV2	164296	272193	9.66%	0.639%
30	9.109	2285	2292	2305	rVB5	42648	85004	3.02%	0.200%
31	9.235	2321	2331	2345	rVB	73872	143249	5.09%	0.336%
32	9.331	2346	2361	2373	rBV3	40646	99944	3.55%	0.235%
33	9.402	2373	2383	2386	rBV2	38519	62636	2.22%	0.147%
34	9.486	2399	2409	2428	rVB3	122185	253659	9.00%	0.596%

LSC Area Percent Report

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Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-2

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

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Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N112320W.M
Title : SW846 8260

35	9.582	2429	2439	2445	rBV2	104023	200231	7.11%	0.470%
36	9.624	2446	2452	2466	rVB4	135854	280160	9.95%	0.658%
37	9.836	2498	2518	2535	rBV7	61381	236312	8.39%	0.555%
38	9.936	2539	2549	2558	rVV2	87442	158010	5.61%	0.371%
39	10.058	2574	2587	2602	rBV2	645301	1272910	45.19%	2.989%
40	10.232	2632	2641	2644	rBV5	47855	79261	2.81%	0.186%
41	10.354	2674	2679	2686	rVB4	23190	32215	1.14%	0.076%
42	10.402	2686	2694	2702	rBV5	31443	51108	1.81%	0.120%
43	10.489	2710	2721	2738	rVB5	182053	390034	13.85%	0.916%
44	10.595	2742	2754	2765	rVB9	27864	75636	2.69%	0.178%
45	10.691	2765	2784	2791	rBV5	17599	51027	1.81%	0.120%
46	10.762	2791	2806	2813	rBV7	42327	107729	3.82%	0.253%
47	10.891	2839	2846	2851	rBV4	37463	52641	1.87%	0.124%
48	11.035	2883	2891	2899	rBV7	20340	37823	1.34%	0.089%
49	11.087	2900	2907	2917	rVB6	26523	47370	1.68%	0.111%
50	11.180	2919	2936	2954	rVB10	40260	127446	4.52%	0.299%
51	11.273	2954	2965	2981	rBV5	33315	98903	3.51%	0.232%
52	11.379	2986	2998	3011	rBV	578935	1024547	36.37%	2.406%
53	11.482	3019	3030	3048	rVB	658637	1152662	40.92%	2.707%
54	11.598	3049	3066	3085	rVV	717512	1342916	47.67%	3.154%
55	11.724	3098	3105	3116	rVB9	19970	33591	1.19%	0.079%
56	11.920	3147	3166	3183	rVB3	39835	109694	3.89%	0.258%
57	12.222	3249	3260	3272	rBV	360814	584842	20.76%	1.373%
58	12.373	3295	3307	3319	rBV	465164	772890	27.44%	1.815%
59	12.563	3354	3366	3379	rVV	617226	986289	35.01%	2.316%
60	12.659	3379	3396	3404	rVV	234196	415928	14.77%	0.977%
61	12.707	3404	3411	3421	rVB	104100	169159	6.01%	0.397%
62	12.865	3447	3460	3474	rBV	462850	760382	26.99%	1.786%
63	13.013	3494	3506	3519	rVB	1115500	1752072	62.20%	4.114%
64	13.141	3531	3546	3558	rBV3	67057	164672	5.85%	0.387%
65	13.264	3575	3584	3589	rBV2	22819	33228	1.18%	0.078%
66	13.315	3589	3600	3605	rBV	598438	1010365	35.87%	2.373%
67	13.347	3605	3610	3622	rVB	692699	1062185	37.71%	2.494%
68	13.486	3641	3653	3654	rBV2	441581	490293	17.41%	1.151%
69	13.514	3655	3662	3672	rVV	1689384	2816940	100.00%	6.615%
70	13.566	3672	3678	3693	rVV3	190036	406837	14.44%	0.955%
71	13.646	3694	3703	3711	rVB2	64208	99061	3.52%	0.233%

LSC Area Percent Report

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Sample : L4829-12
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-2

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

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Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N112320W.M
Title : SW846 8260

72	13.711	3711	3723	3734	rVB	171238	282220	10.02%	0.663%
73	13.778	3734	3744	3759	rBV3	171873	381272	13.53%	0.895%
74	13.862	3759	3770	3780	rVV	791572	1211099	42.99%	2.844%
75	13.920	3780	3788	3794	rVV	168376	268714	9.54%	0.631%
76	13.968	3794	3803	3816	rVV	890637	1429350	50.74%	3.356%
77	14.039	3818	3825	3835	rVB6	34847	59636	2.12%	0.140%
78	14.100	3835	3844	3853	rVB2	84391	130611	4.64%	0.307%
79	14.180	3853	3869	3876	rBV	600559	978104	34.72%	2.297%
80	14.219	3876	3881	3890	rVV	307097	472576	16.78%	1.110%
81	14.267	3890	3896	3912	rVB4	75483	140087	4.97%	0.329%
82	14.402	3931	3938	3943	rBV2	37962	51042	1.81%	0.120%
83	14.447	3943	3952	3962	rVV	538205	862593	30.62%	2.026%
84	14.495	3963	3967	3975	rVB2	62138	83358	2.96%	0.196%
85	14.563	3975	3988	4000	rBV3	896361	1793437	63.67%	4.211%
86	14.627	4000	4008	4019	rVB4	130719	229517	8.15%	0.539%
87	14.711	4025	4034	4044	rBV	128469	207779	7.38%	0.488%
88	14.820	4050	4068	4076	rBV5	272209	643438	22.84%	1.511%
89	14.929	4089	4102	4116	rVB2	248127	553035	19.63%	1.299%
90	15.100	4136	4155	4168	rBV	593047	1068174	37.92%	2.508%
91	15.206	4178	4188	4197	rBV4	79818	153138	5.44%	0.360%
92	15.257	4197	4204	4225	rVB	61975	124770	4.43%	0.293%
93	15.383	4231	4243	4256	rBV	103459	191692	6.80%	0.450%
94	15.540	4276	4292	4310	rBV4	58626	200453	7.12%	0.471%
95	15.662	4320	4330	4340	rBV2	40372	83525	2.97%	0.196%
96	15.730	4343	4351	4380	rVB5	61747	157282	5.58%	0.369%
97	15.871	4382	4395	4403	rBV6	18581	38880	1.38%	0.091%
98	16.058	4434	4453	4461	rBV6	17071	43864	1.56%	0.103%
99	16.119	4461	4472	4487	rVB2	82566	164645	5.84%	0.387%
100	16.305	4516	4530	4550	rBV	219779	478296	16.98%	1.123%

Sum of corrected areas: 42584874

LSC Report - Integrated Chromatogram

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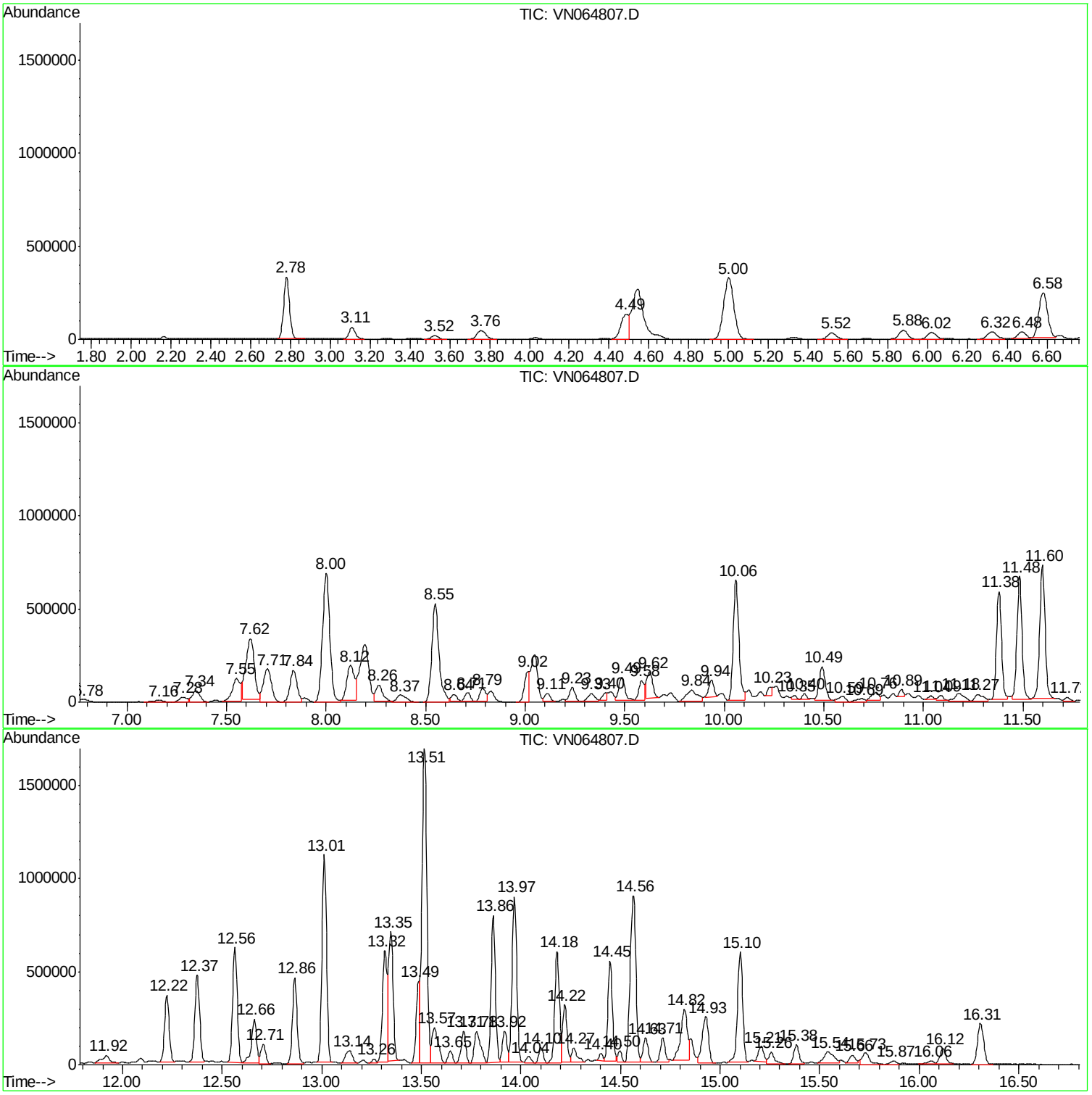
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N112320W.M
Quant Title : SW846 8260

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



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Quant Title : SW846 8260

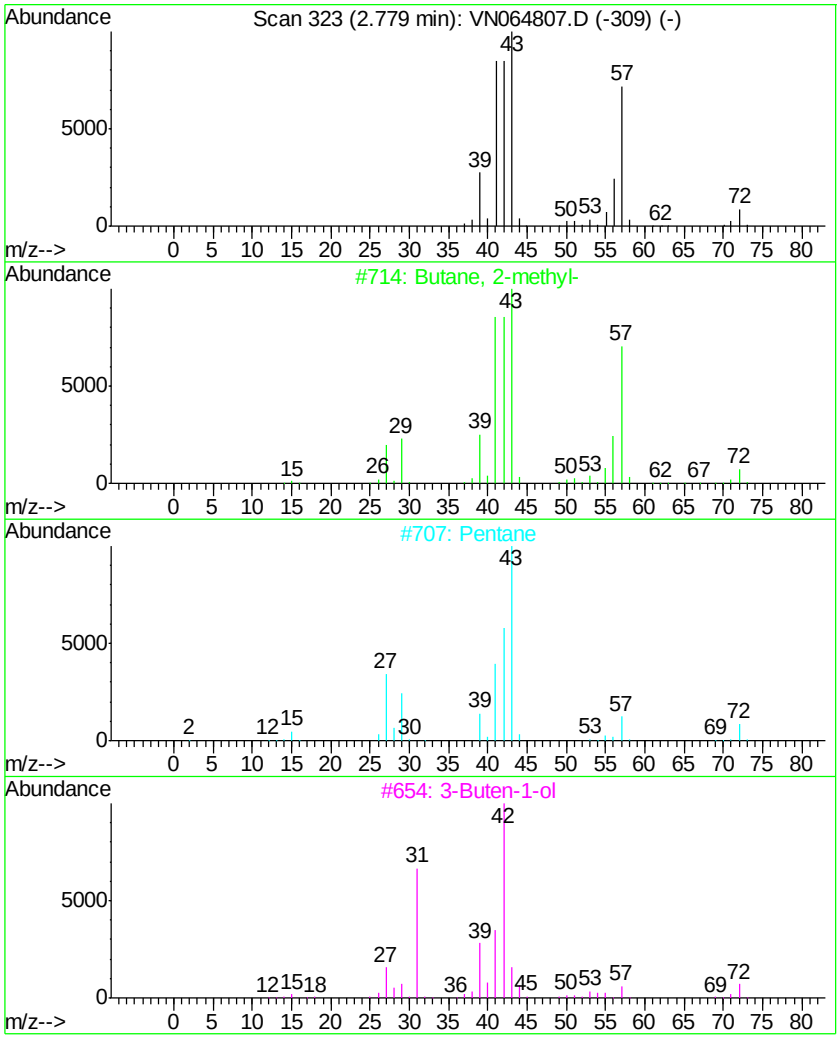
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

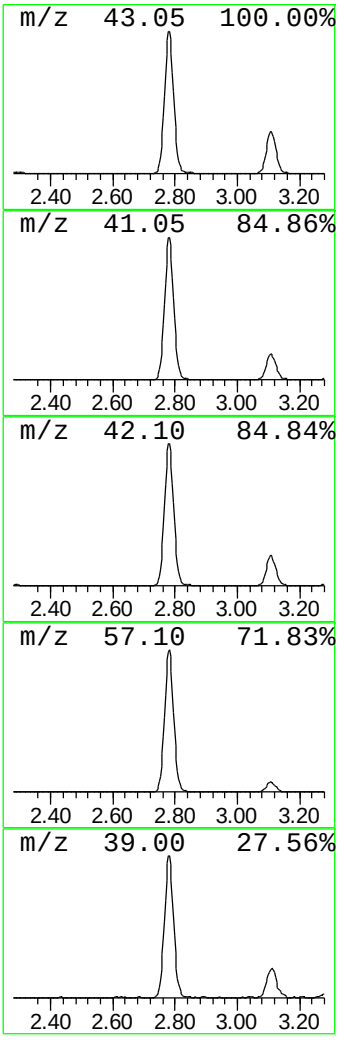
****

Peak Number 1 Butane, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.78	34.43 ug/l	670592	Pentafluorobenzene	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	94
2	Pentane	72	C5H12	000109-66-0	45
3	3-Buten-1-ol	72	C4H8O	000627-27-0	28
4	1-Propene, 2-methyl-	56	C4H8	000115-11-7	9
5	Methane, isocyanato-	57	C2H3NO	000624-83-9	9





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ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N112320W.M
Quant Title : SW846 8260

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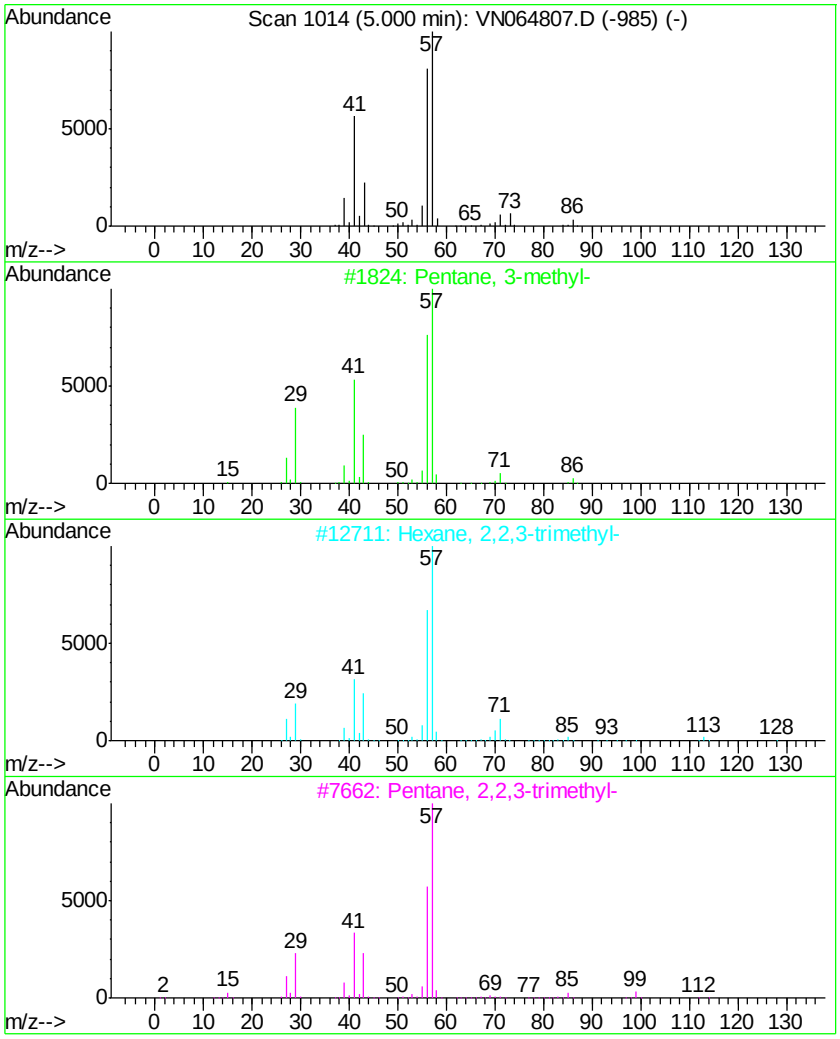
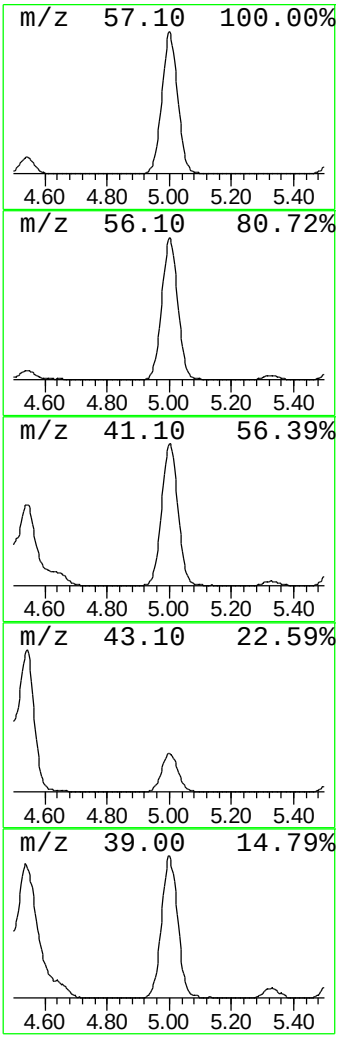
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.00	61.29 ug/l	1193670	Pentafluorobenzene	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	78
3	Pentane, 2,2,3-trimethyl-	114	C8H18	000564-02-3	50
4	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	45
5	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	42





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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
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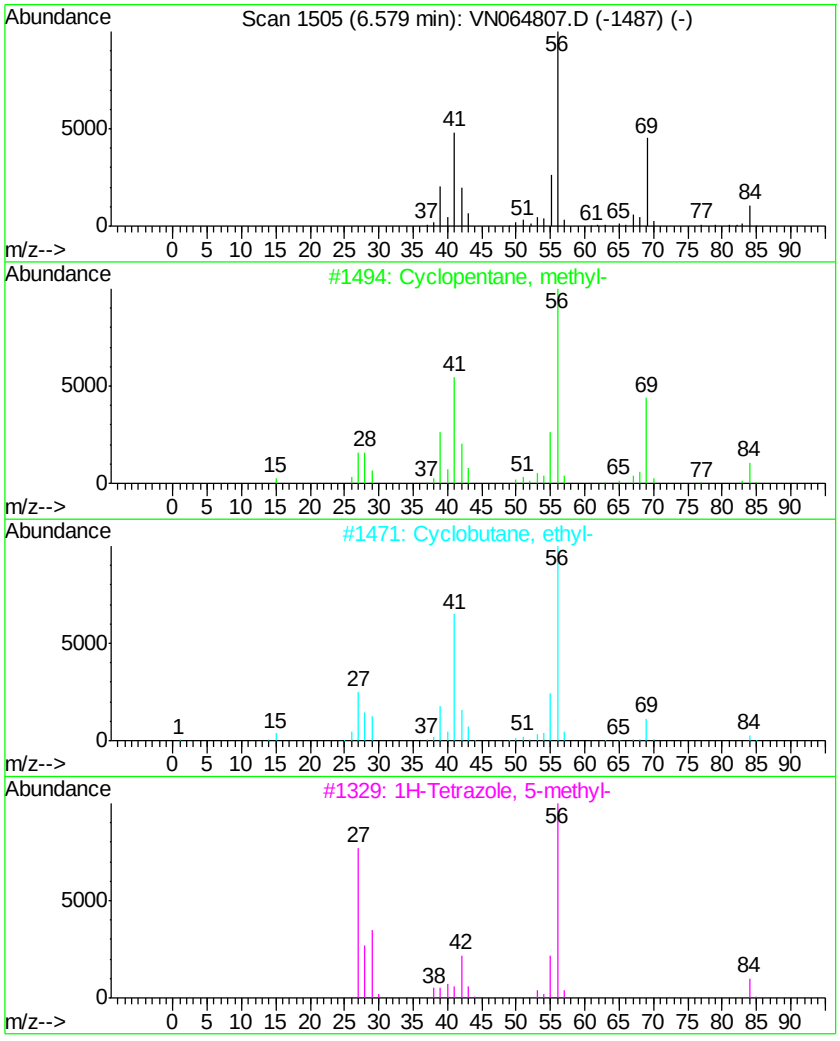
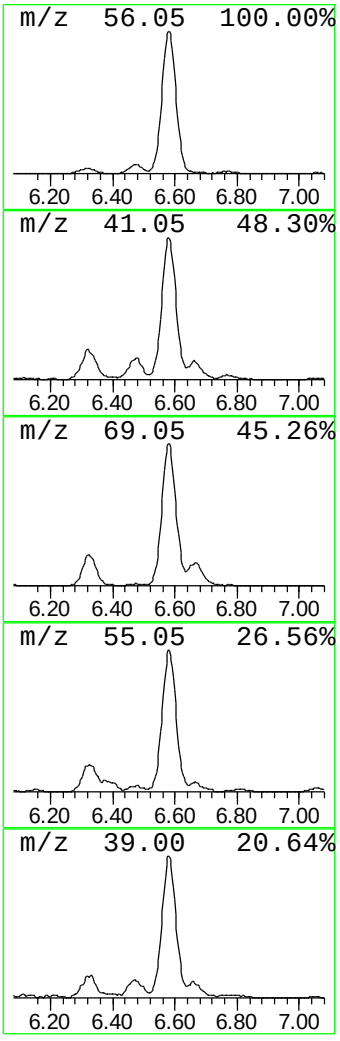
Peak Number 3 Cyclopentane, methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	37.90 ug/l	738021	Pentafluorobenzene	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
3	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64
4	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	43
5	2H-Pyran-2,6(3H)-dione, dihydro-...	142	C7H10O3	004160-82-1	38





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Instrument : MSVOA_N
ClientSampleId : MW-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N112320W.M
Quant Title : SW846 8260

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

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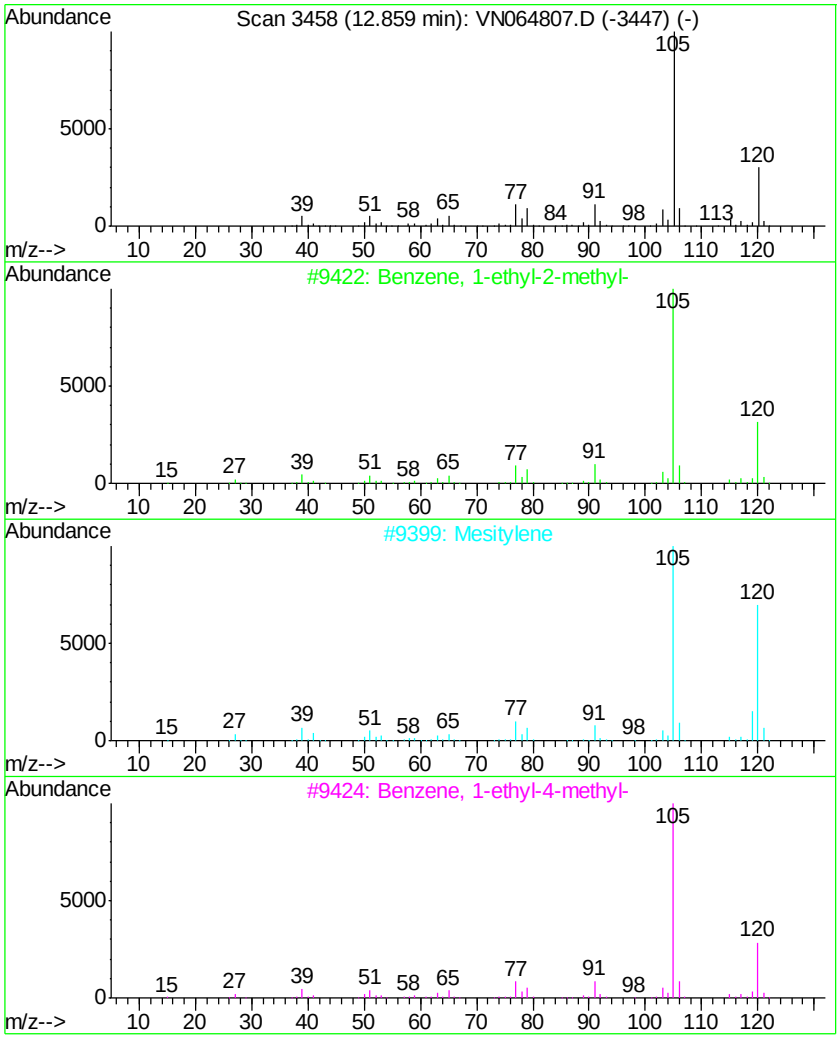
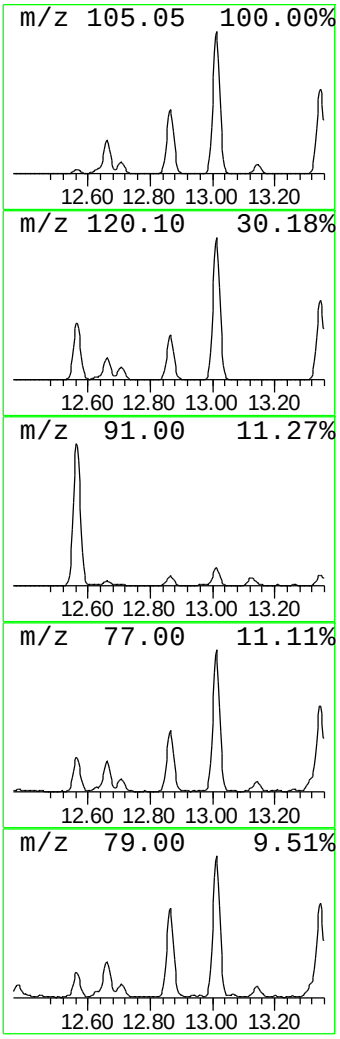
Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.86	37.63 ug/l	760382	1,4-Dichlorobenzene-d4	13.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2	Mesitylene	120	C9H12	000108-67-8	91
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91





Library Search Compound Report

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Sample : L4829-12
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N112320W.M
Quant Title : SW846 8260

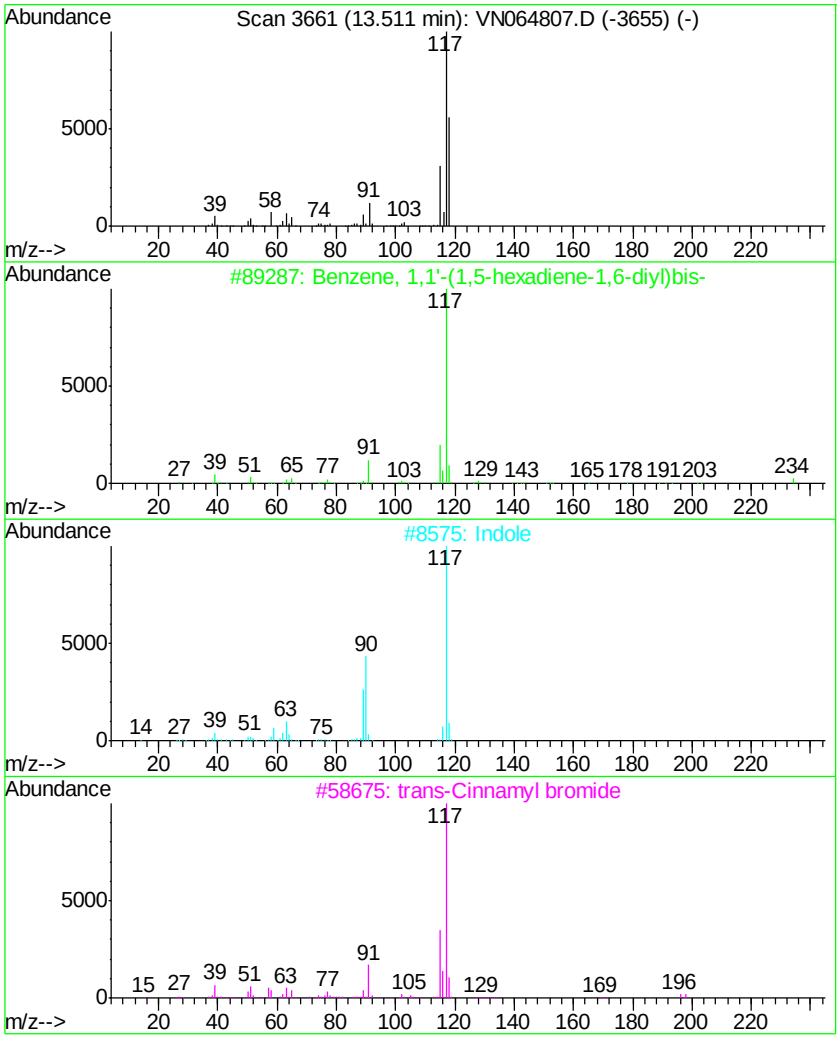
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

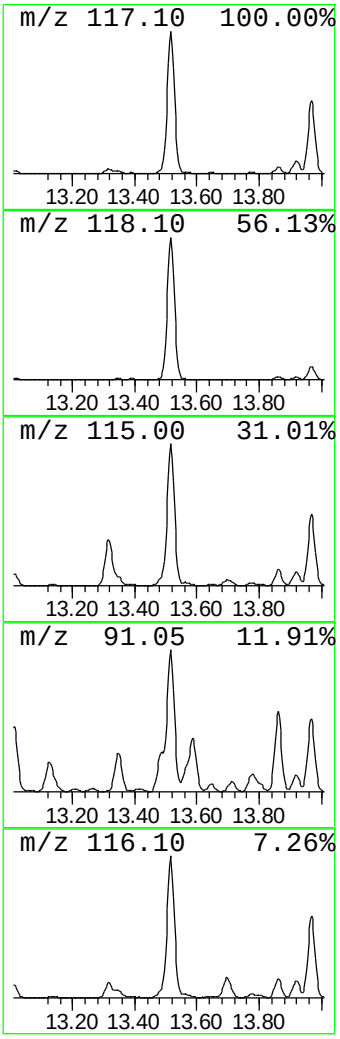
****

Peak Number 5 Benzene, 1,1'-(1,5-hexadien... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.51	139.40 ug/l	2816940	1,4-Dichlorobenzene-d4	13.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	53
2	Indole	117	C8H7N	000120-72-9	46
3	trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	42
4	Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	40
5	m-Aminophenylacetylene	117	C8H7N	054060-30-9	37





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN112420\
Data File : VN064807.D
Acq On : 24 Nov 2020 13:50
Operator : JC/MD
Sample : L4829-12
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N112320W.M
Quant Title : SW846 8260

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

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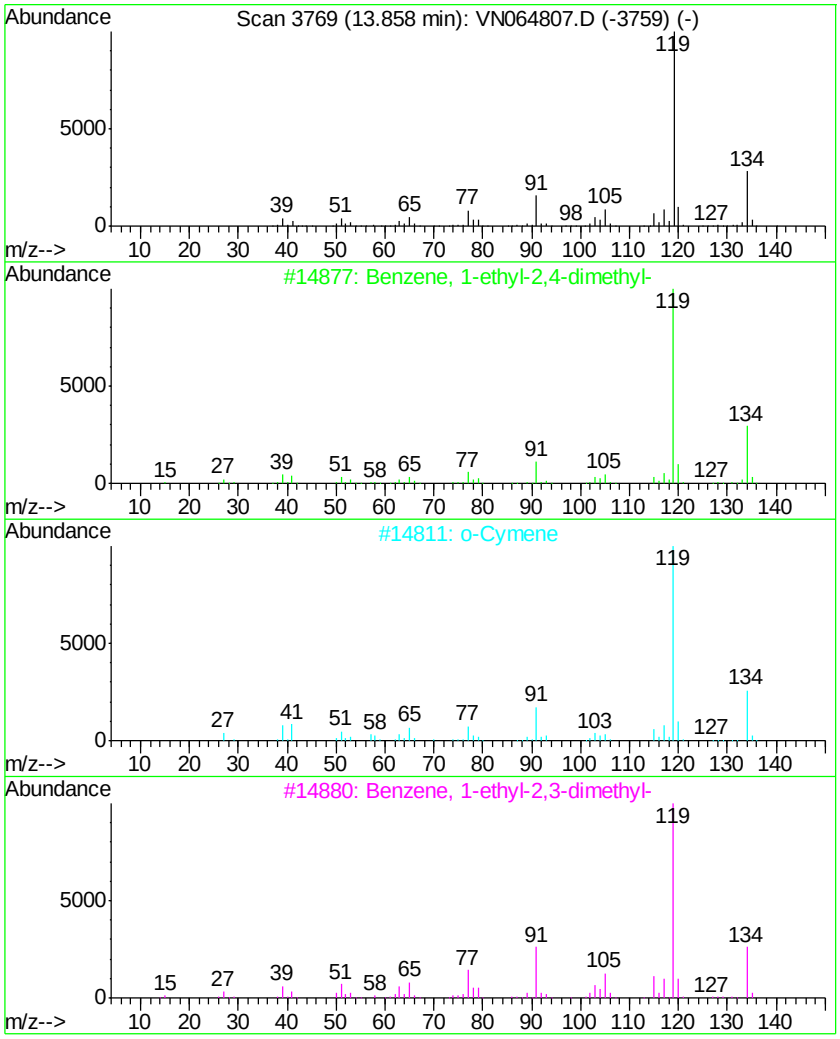
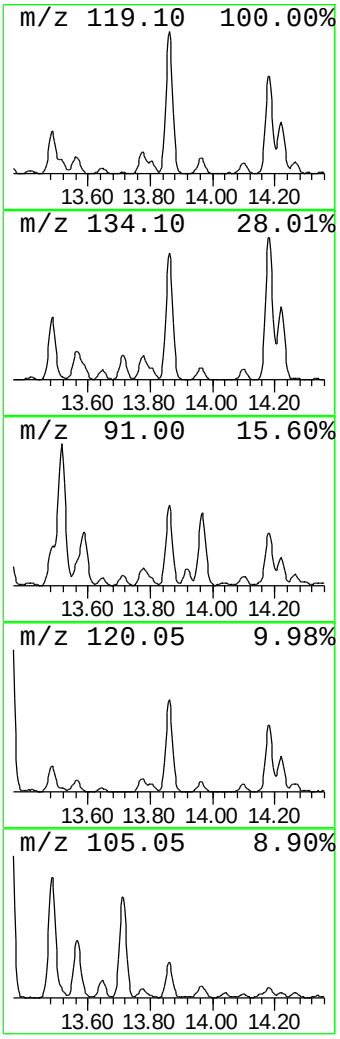
Peak Number 6 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	59.93 ug/l	1211100	1,4-Dichlorobenzene-d4	13.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2	o-Cymene	134	C10H14	000527-84-4	97
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN112420\
Data File : VN064807.D
Acq On : 24 Nov 2020 13:50
Operator : JC/MD
Sample : L4829-12
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N112320W.M
Quant Title : SW846 8260

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

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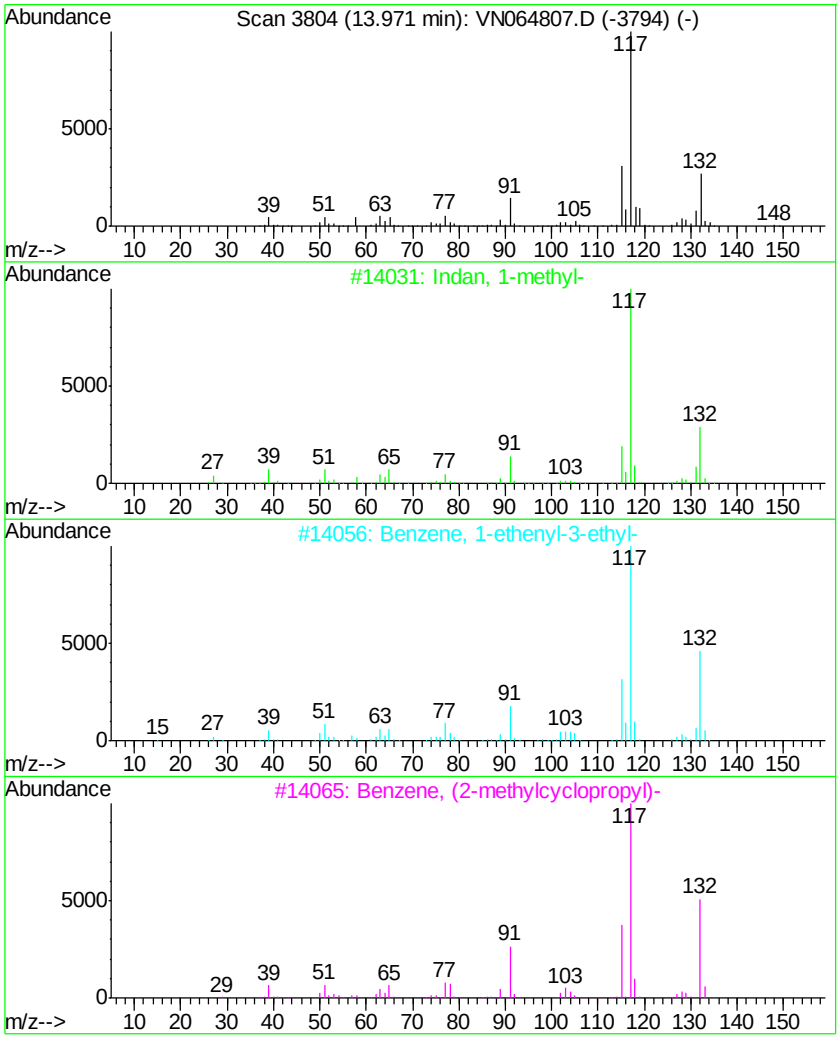
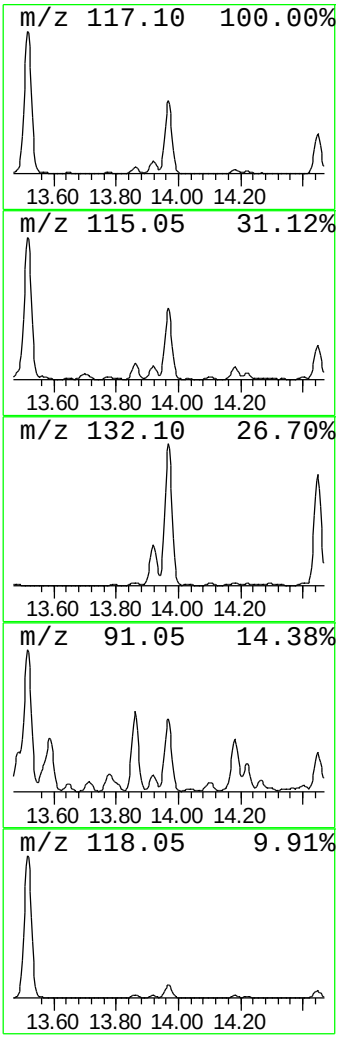
Peak Number 7 Indan, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	70.73 ug/l	1429350	1,4-Dichlorobenzene-d4	13.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	93
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
3	Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	81
4	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	80
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	80





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN112420\
Data File : VN064807.D
Acq On : 24 Nov 2020 13:50
Operator : JC/MD
Sample : L4829-12
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N112320W.M
Quant Title : SW846 8260

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

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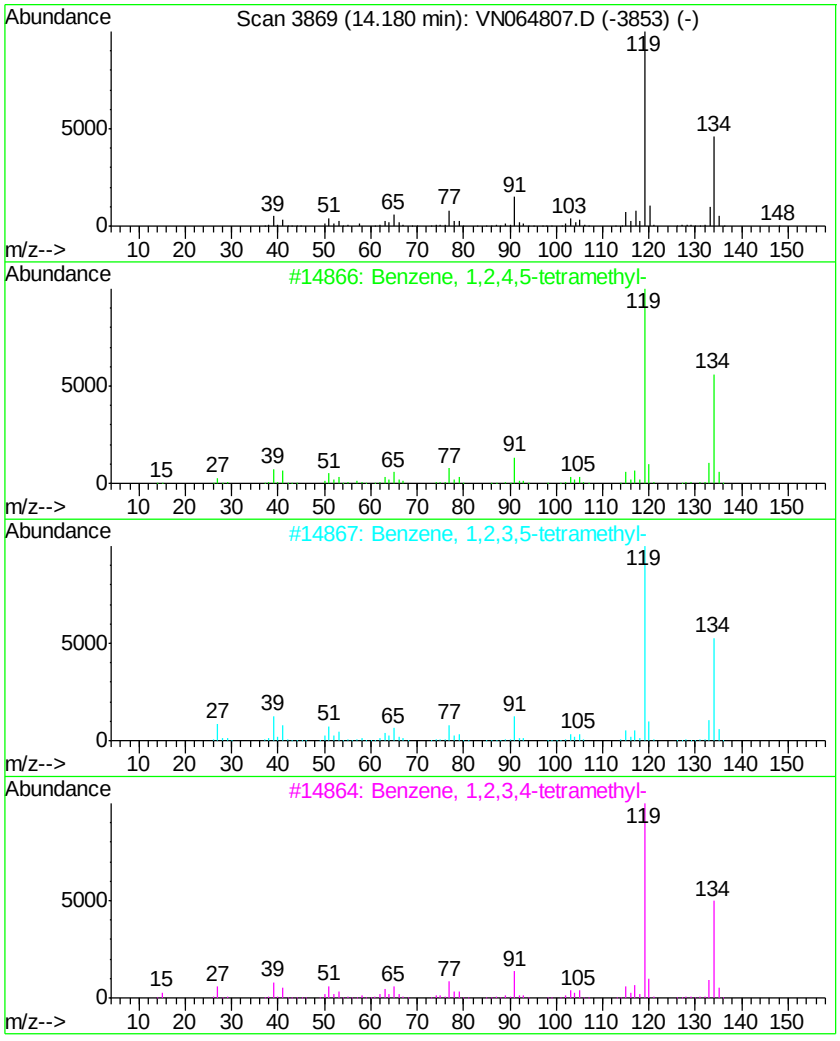
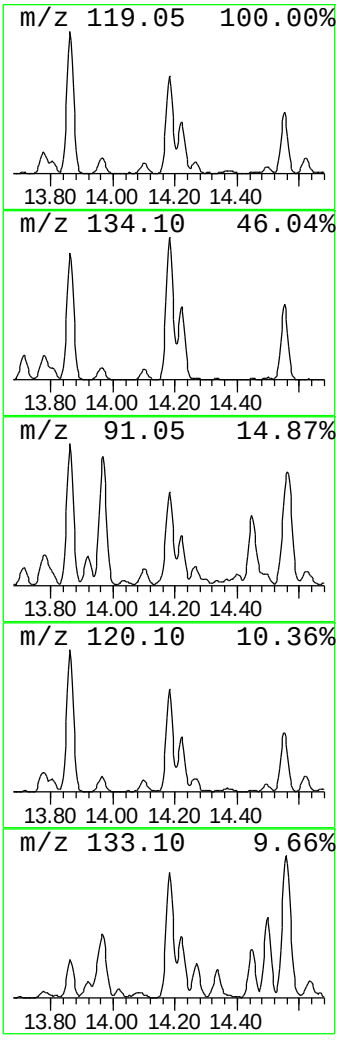
Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.18	48.40 ug/l	978104	1,4-Dichlorobenzene-d4	13.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94





Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN112420\
Data File : VN064807.D
Acq On : 24 Nov 2020 13:50
Operator : JC/MD
Sample : L4829-12
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N112320W.M
Quant Title : SW846 8260

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

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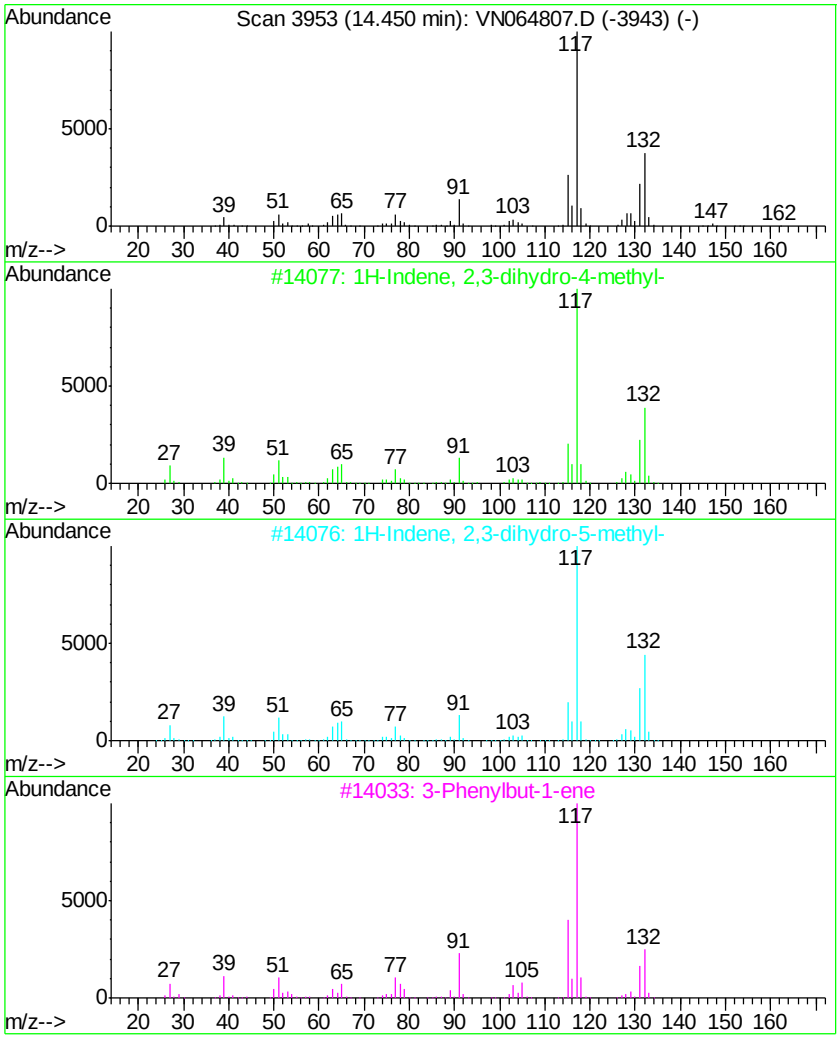
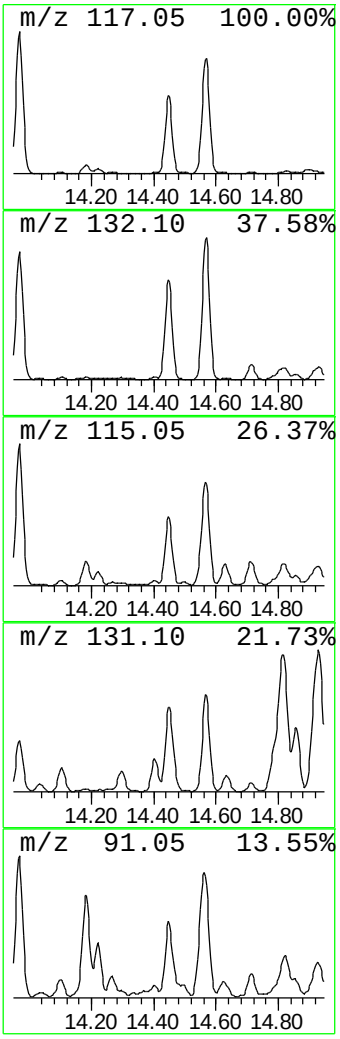
Peak Number 9 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.45	42.69 ug/l	862593	1,4-Dichlorobenzene-d4	13.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	95
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
4	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	87
5	1-Phenyl-1-butene	132	C10H12	000824-90-8	86





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN112420\
Data File : VN064807.D
Acq On : 24 Nov 2020 13:50
Operator : JC/MD
Sample : L4829-12
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N112320W.M
Quant Title : SW846 8260

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

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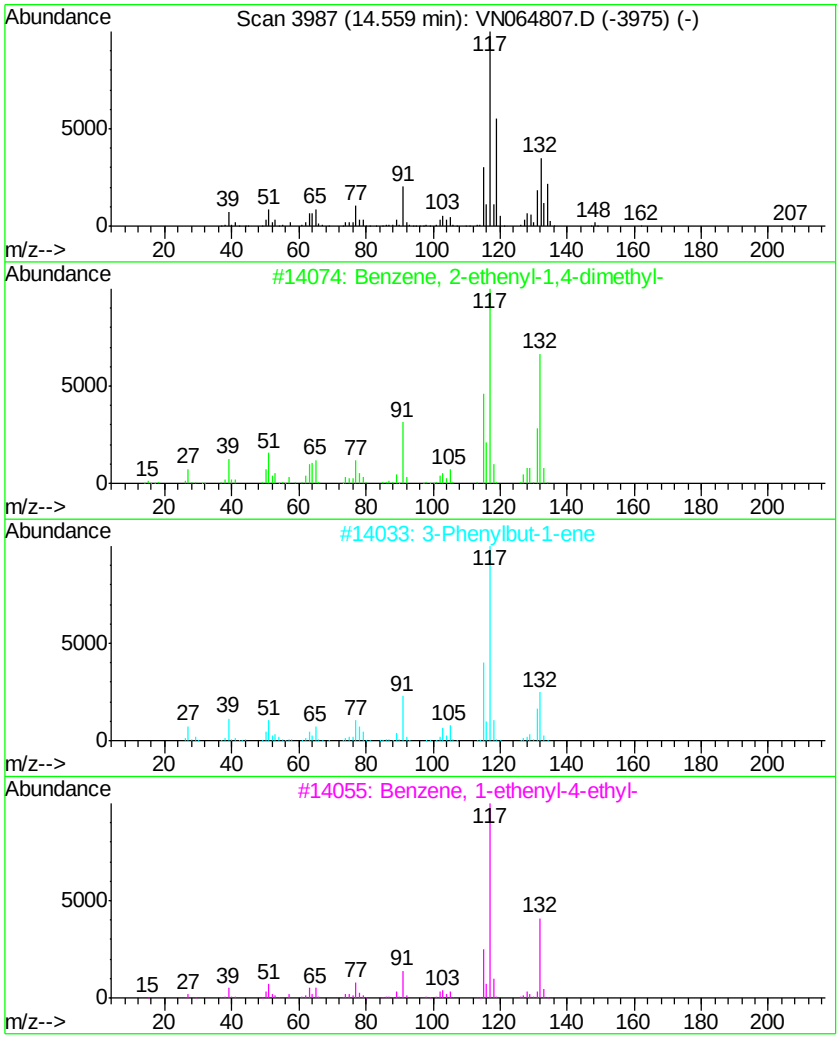
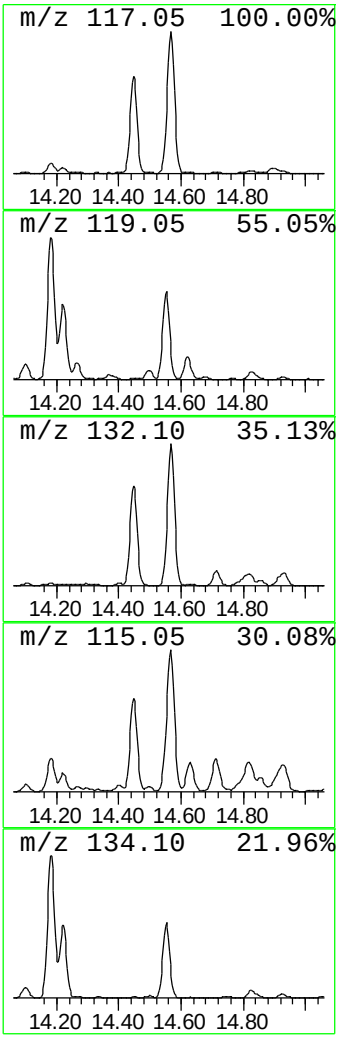
Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.56	88.75 ug/l	1793440	1,4-Dichlorobenzene-d4	13.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
3	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	70
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	68
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	68





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN112420\
Data File : VN064807.D
Acq On : 24 Nov 2020 13:50
Operator : JC/MD
Sample : L4829-12
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N112320W.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	34.4	ug/l	670592	1	7.63	973713	50.0
Pentane, 3-methyl-	5.00	61.3	ug/l	1193670	1	7.63	973713	50.0
Cyclopentane, met...	6.58	37.9	ug/l	738021	1	7.63	973713	50.0
Benzene, 1-ethyl-...	12.86	37.6	ug/l	760382	4	13.32	1010370	50.0
Benzene, 1,1'-(1,...	13.51	139.4	ug/l	2816940	4	13.32	1010370	50.0
Benzene, 1-ethyl-...	13.86	59.9	ug/l	1211100	4	13.32	1010370	50.0
Indan, 1-methyl-	13.97	70.7	ug/l	1429350	4	13.32	1010370	50.0
Benzene, 1,2,4,5-...	14.18	48.4	ug/l	978104	4	13.32	1010370	50.0
1H-Indene, 2,3-di...	14.45	42.7	ug/l	862593	4	13.32	1010370	50.0
Benzene, 2-etheny...	14.56	88.8	ug/l	1793440	4	13.32	1010370	50.0