LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN022020\
Data File : VN060167.D
Acq On : 20 Feb 2020 12:53
Operator : JC/MD
Sample : L1589-01 10X
Misc : 11.95g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : SB3-9.5-10.0

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\82N021220W.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	7.564	1784	1801	1811	rBV	116686	296806	4.75%	0.288%
2	7.641	1812	1825	1841	rVB2	264925	670030	10.71%	0.651%
3	7.854	1877	1891	1904	rBV2	30883	74202	1.19%	0.072%
4	8.005	1920	1938	1963	rBV2	146272	368701	5.89%	0.358%
5	8.416	2052	2066	2082	rBV	86315	178059	2.85%	0.173%
6	8.567	2094	2113	2136	rBV	320721	671787	10.74%	0.653%
7	9.063	2242	2267	2281	rBV2	291065	689343	11.02%	0.670%
8	9.255	2313	2327	2339	rVV	51842	103959	1.66%	0.101%
9	9.497	2388	2402	2415	rBV4	33075	69351	1.11%	0.067%
10	9.715	2458	2470	2494	rVB2	120509	291319	4.66%	0.283%
11	9.857	2494	2514	2534	rBV2	81649	208530	3.33%	0.203%
12	10.075	2568	2582	2592	rVV3	673488	1299214	20.77%	1.262%
13	10.143	2592	2603	2615	rVV	1019715	1926963	30.81%	1.872%
14	10.204	2615	2622	2627	rVV2	53695	94597	1.51%	0.092%
15	10.271	2628	2643	2658	rVB	381908	763030	12.20%	0.741%
16	10.419	2671	2689	2705	rVB	92450	194492	3.11%	0.189%
17	10.519	2705	2720	2732	rBV2	109349	226213	3.62%	0.220%
18	10.709	2769	2779	2787	rBV	74708	119779	1.92%	0.116%
19	10.873	2821	2830	2836	rBV2	71118	118960	1.90%	0.116%
20	10.940	2844	2851	2860	rVB	183530	286936	4.59%	0.279%
21	10.995	2860	2868	2880	rVB2	141405	252453	4.04%	0.245%
22	11.114	2892	2905	2913	rBV5	64640	130391	2.08%	0.127%
23	11.194	2913	2930	2949	rVB6	199992	478869	7.66%	0.465%
24	11.291	2949	2960	2971	rBV	171291	305232	4.88%	0.297%
25	11.397	2982	2993	3012	rVV	483585	846898	13.54%	0.823%
26	11.503	3012	3026	3042	rVV	1130633	2006913	32.09%	1.950%
27	11.612	3042	3060	3078	rVV2	2859570	6254685	100.00%	6.078%
28	11.689	3079	3084	3095	rVB4	111643	183983	2.94%	0.179%
29	11.940	3140	3162	3178	rBV	1528387	2909410	46.52%	2.827%
30	12.040	3185	3193	3200	rBV	113784	158278	2.53%	0.154%
31	12.114	3209	3216	3221	rBV3	88882	140967	2.25%	0.137%
32	12.159	3222	3230	3246	rVV7	215321	496215	7.93%	0.482%
33	12.242	3247	3256	3267	rVB	617610	983148	15.72%	0.955%
34	12.397	3296	3304	3313	rVB	408372	627668	10.04%	0.610%

LSC Area Percent Report

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Sample : L1589-01 10X
Misc : 11.95g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : SB3-9.5-10.0

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\82N021220W.M
Title : SW846 8260

35	12.442	3313	3318	3338	rVB4	90815	168610	2.70%	0.164%
36	12.586	3347	3363	3372	rBV	883234	1487474	23.78%	1.445%
37	12.651	3372	3383	3389	rBV	2388923	3851563	61.58%	3.743%
38	12.725	3399	3406	3417	rVB3	1251220	2000924	31.99%	1.944%
39	12.773	3417	3421	3431	rVB2	84505	121787	1.95%	0.118%
40	12.886	3443	3456	3470	rBV	1417715	2364933	37.81%	2.298%
41	12.956	3471	3478	3490	rVB3	196813	335870	5.37%	0.326%
42	13.033	3490	3502	3514	rBV	2670492	4266079	68.21%	4.145%
43	13.165	3530	3543	3552	rBV	773955	1328885	21.25%	1.291%
44	13.230	3552	3563	3572	rVV	1008886	1613635	25.80%	1.568%
45	13.284	3572	3580	3587	rVV	406336	599976	9.59%	0.583%
46	13.342	3588	3598	3601	rVV	618947	1012202	16.18%	0.984%
47	13.368	3602	3606	3616	rVB	958183	1362020	21.78%	1.324%
48	13.422	3617	3623	3640	rBV3	63494	180246	2.88%	0.175%
49	13.538	3641	3659	3666	rBV	1897438	3654694	58.43%	3.551%
50	13.583	3666	3673	3689	rVB4	1304881	2752241	44.00%	2.674%
51	13.667	3689	3699	3708	rBV3	800482	1264848	20.22%	1.229%
52	13.734	3711	3720	3730	rVB	815189	1245737	19.92%	1.211%
53	13.799	3731	3740	3744	rBV	822904	1218312	19.48%	1.184%
54	13.828	3745	3749	3758	rVB	760487	1018203	16.28%	0.989%
55	13.885	3758	3767	3776	rVB2	1046955	1579821	25.26%	1.535%
56	13.940	3776	3784	3789	rBV	194643	284330	4.55%	0.276%
57	13.992	3790	3800	3810	rVB2	1686571	2787529	44.57%	2.709%
58	14.046	3810	3817	3819	rBV	168805	205632	3.29%	0.200%
59	14.072	3821	3825	3832	rVV2	237395	303543	4.85%	0.295%
60	14.123	3833	3841	3847	rVV2	309367	473704	7.57%	0.460%
61	14.165	3848	3854	3860	rVV6	197075	324583	5.19%	0.315%
62	14.207	3861	3867	3872	rVV2	407982	621414	9.94%	0.604%
63	14.242	3872	3878	3886	rVB	747123	1065938	17.04%	1.036%
64	14.291	3886	3893	3900	rBV2	604079	847697	13.55%	0.824%
65	14.329	3900	3905	3912	rVB	331292	417967	6.68%	0.406%
66	14.429	3931	3936	3942	rVB2	406368	504701	8.07%	0.490%
67	14.471	3942	3949	3958	rBV3	891004	1519636	24.30%	1.477%
68	14.519	3959	3964	3973	rVB	1023679	1462356	23.38%	1.421%
69	14.586	3973	3985	3997	rBV5	2023803	4593117	73.43%	4.463%
70	14.648	3997	4004	4013	rVB3	377991	599228	9.58%	0.582%
71	14.738	4022	4032	4045	rVB	1530063	2491996	39.84%	2.422%

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ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : SB3-9.5-10.0

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
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Start Thrs: 0.2 Max Peaks: 100
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Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N021220W.M
Title : SW846 8260

72	14.811	4048	4055	4057	rBV4	287259	349511	5.59%	0.340%
73	14.847	4059	4066	4072	rBV4	567055	809427	12.94%	0.787%
74	14.956	4089	4100	4112	rVB2	829015	1743566	27.88%	1.694%
75	15.184	4163	4171	4179	rBV	1025866	1599776	25.58%	1.555%
76	15.249	4180	4191	4208	rVV5	830833	2692435	43.05%	2.616%
77	15.326	4209	4215	4230	rVB7	444599	834457	13.34%	0.811%
78	15.413	4231	4242	4255	rBV3	607562	1241291	19.85%	1.206%
79	15.609	4295	4303	4318	rVB	2292609	4209178	67.30%	4.090%
80	15.696	4324	4330	4340	rVV7	200583	347959	5.56%	0.338%
81	15.763	4342	4351	4374	rVB5	542340	1575052	25.18%	1.531%
82	15.908	4384	4396	4408	rBV	1330909	2641209	42.23%	2.567%
83	16.094	4443	4454	4464	rBV3	1307411	2591446	41.43%	2.518%
84	16.155	4464	4473	4492	rVB2	1401783	2859372	45.72%	2.779%
85	16.348	4521	4533	4544	rBV5	816725	1975359	31.58%	1.920%
86	16.483	4565	4575	4590	rVB3	454510	1081054	17.28%	1.050%

Sum of corrected areas: 102909904

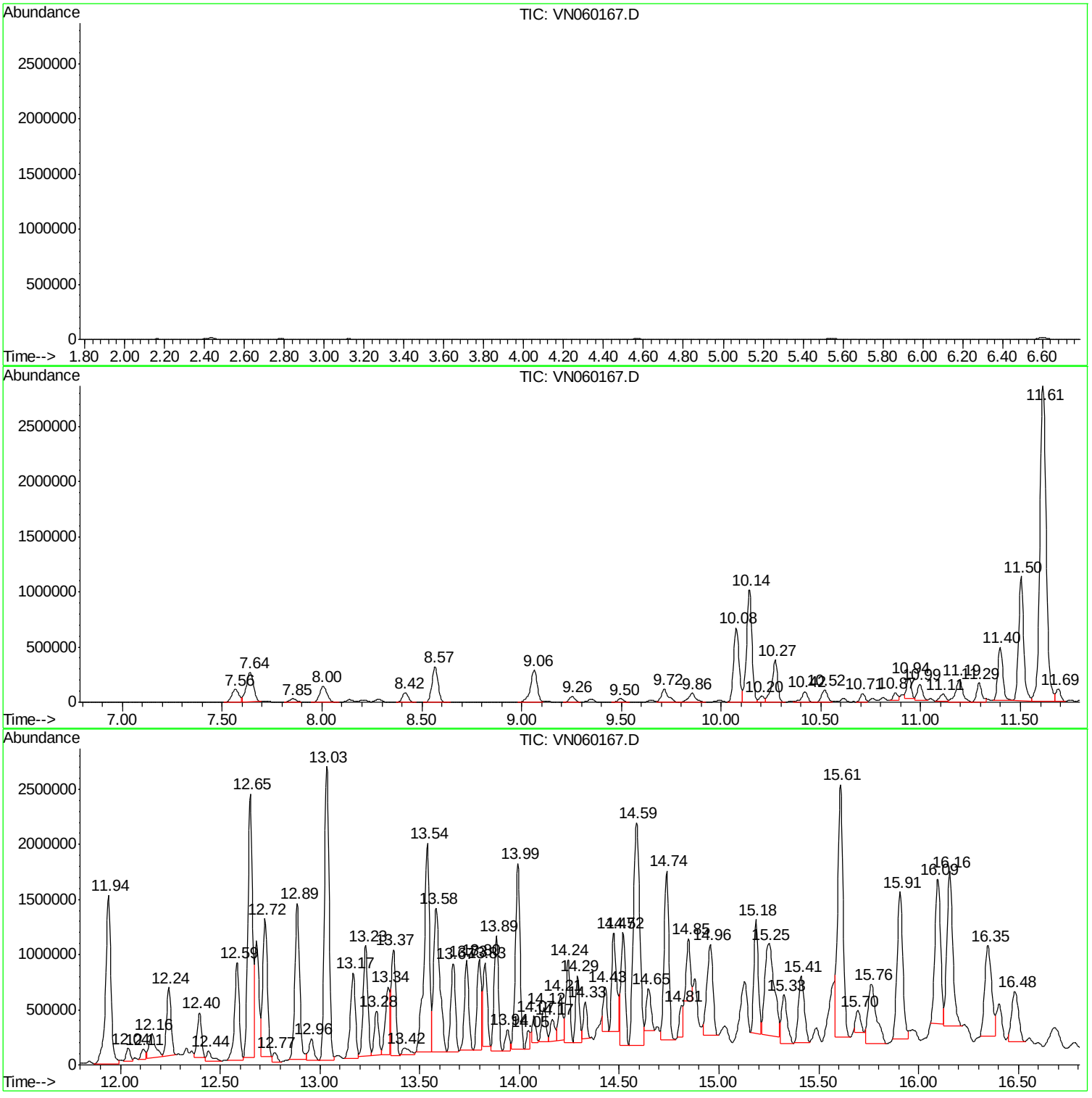
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN022020\
Data File : VN060167.D
Acq On : 20 Feb 2020 12:53
Operator : JC/MD
Sample : L1589-01 10X
Misc : 11.95g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : SB3-9.5-10.0

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N021220W.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\VN022020\
Data File : VN060167.D
Acq On : 20 Feb 2020 12:53
Operator : JC/MD
Sample : L1589-01 10X
Misc : 11.95g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : SB3-9.5-10.0

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N021220W.M
Quant Title : SW846 8260

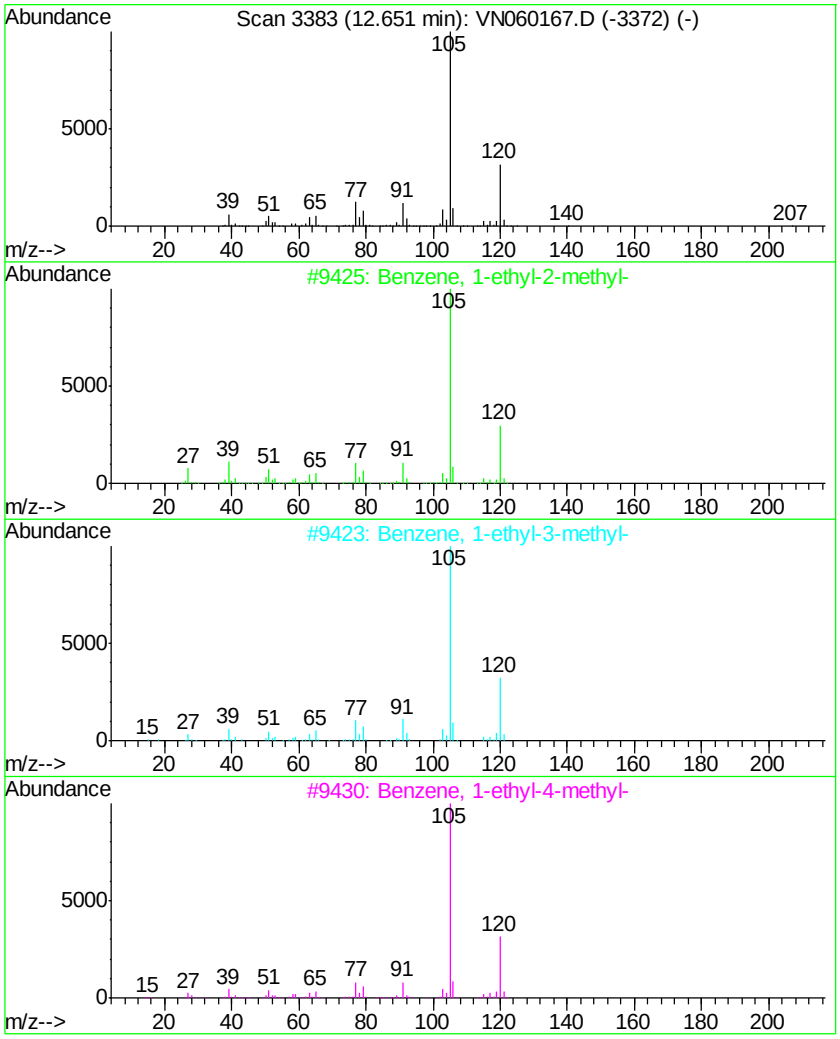
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

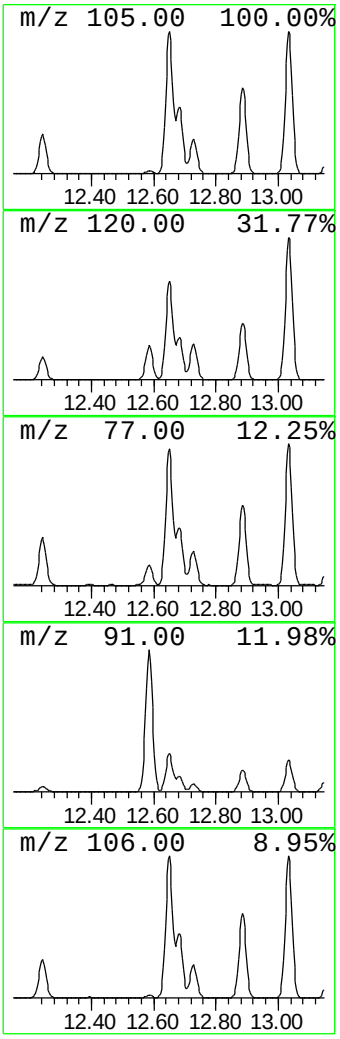
****

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	190.26 ug/l	3851560	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Mesitylene	120	C9H12	000108-67-8	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN022020\
Data File : VN060167.D
Acq On : 20 Feb 2020 12:53
Operator : JC/MD
Sample : L1589-01 10X
Misc : 11.95g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : SB3-9.5-10.0

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N021220W.M
Quant Title : SW846 8260

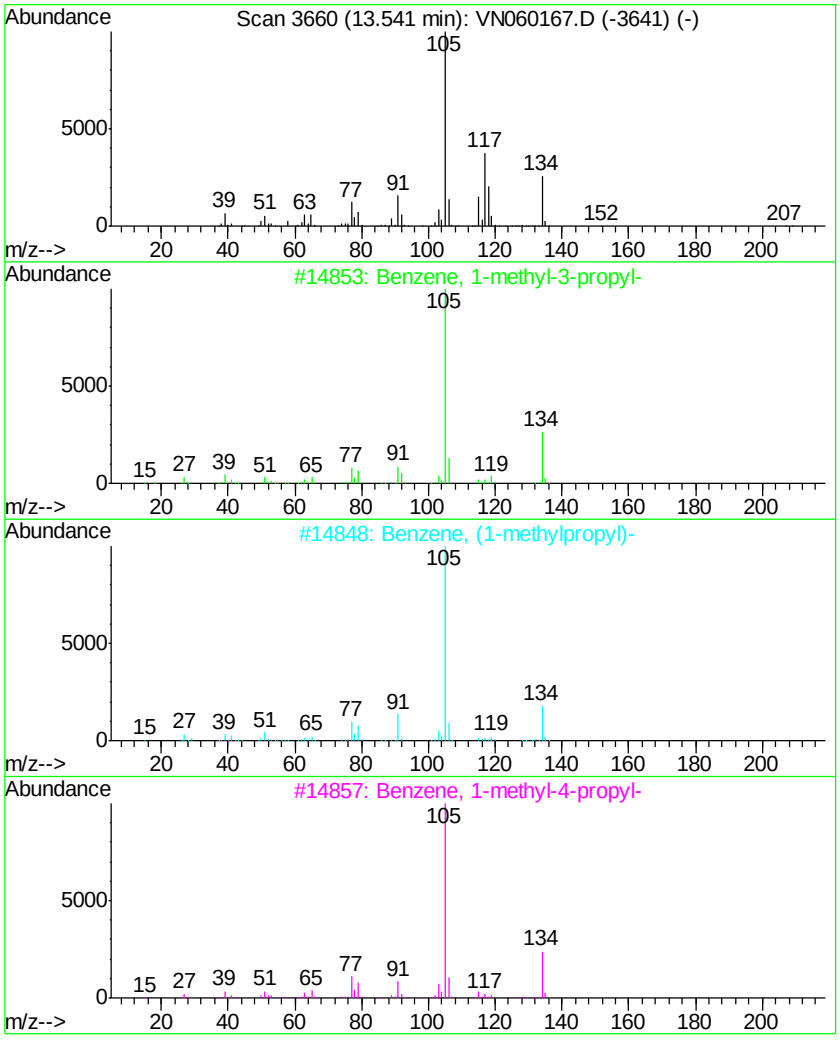
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

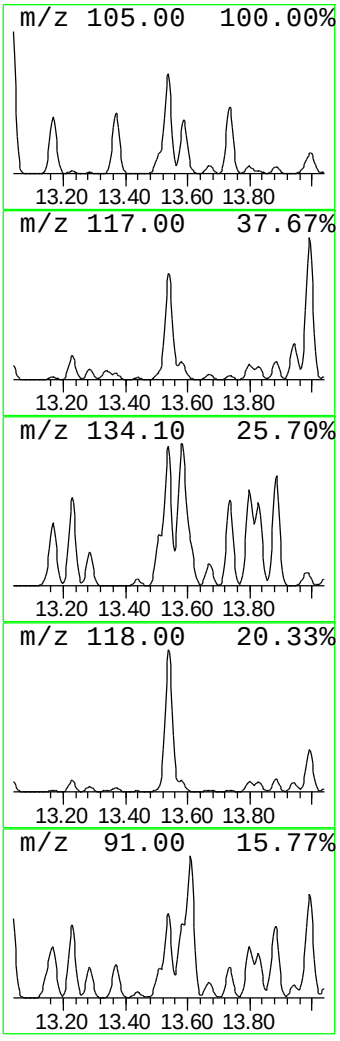
****

Peak Number 2 Benzene, 1-methyl-3-propyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.54	180.53 ug/l	3654690	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	64
2	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	64
3	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	60
4	2-Indanol	134	C9H10O	004254-29-9	58
5	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	49





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ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : SB3-9.5-10.0

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N021220W.M
Quant Title : SW846 8260

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

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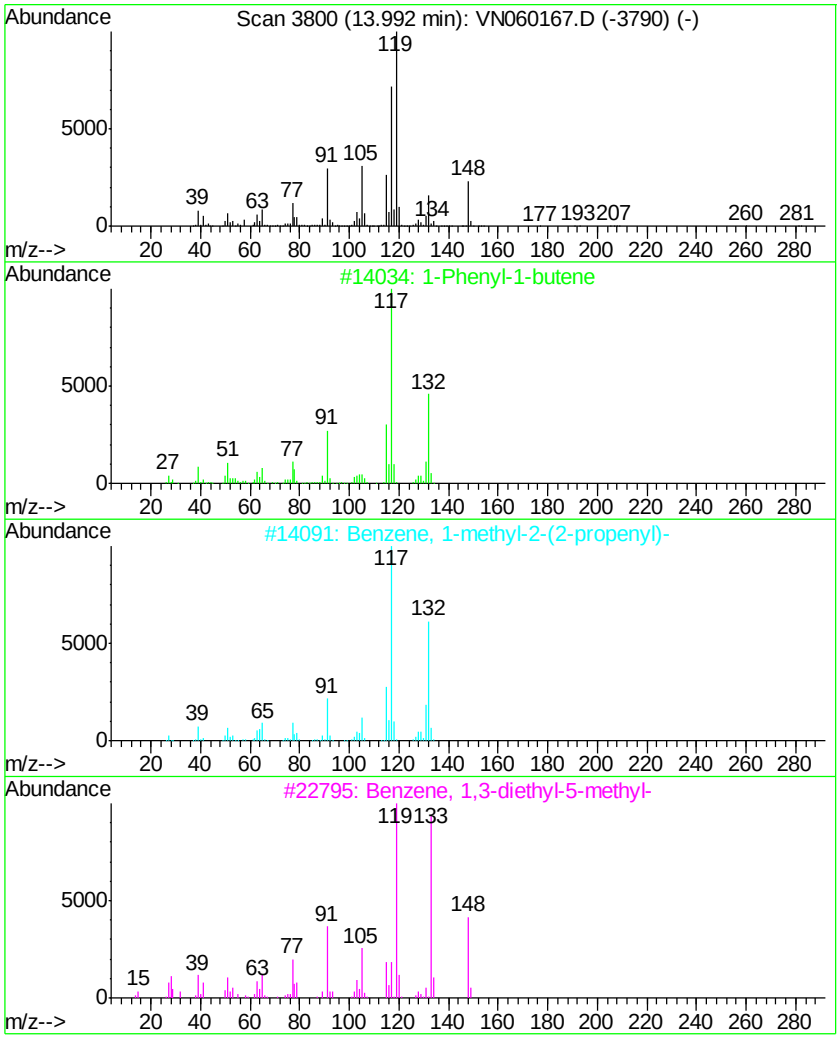
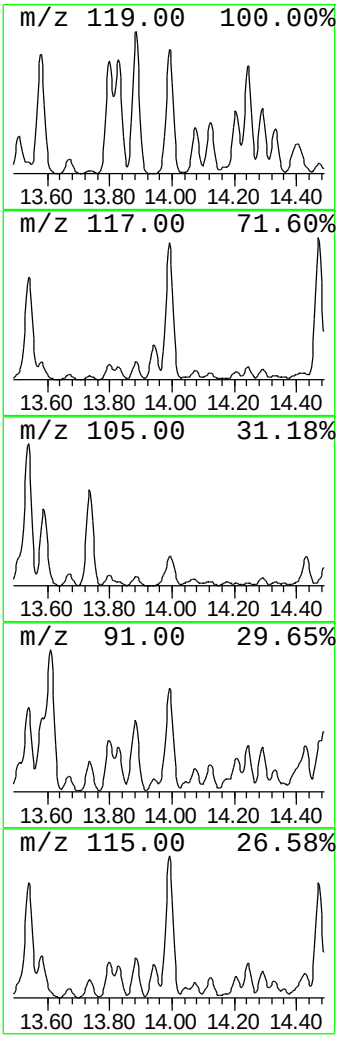
Peak Number 3 unknown13.99 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.99	137.70 ug/l	2787530	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8	49
2	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	46
3	Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	45
4	o-Cymene	134	C10H14	000527-84-4	43
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	42





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN022020\
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Acq On : 20 Feb 2020 12:53
Operator : JC/MD
Sample : L1589-01 10X
Misc : 11.95g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : SB3-9.5-10.0

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N021220W.M
Quant Title : SW846 8260

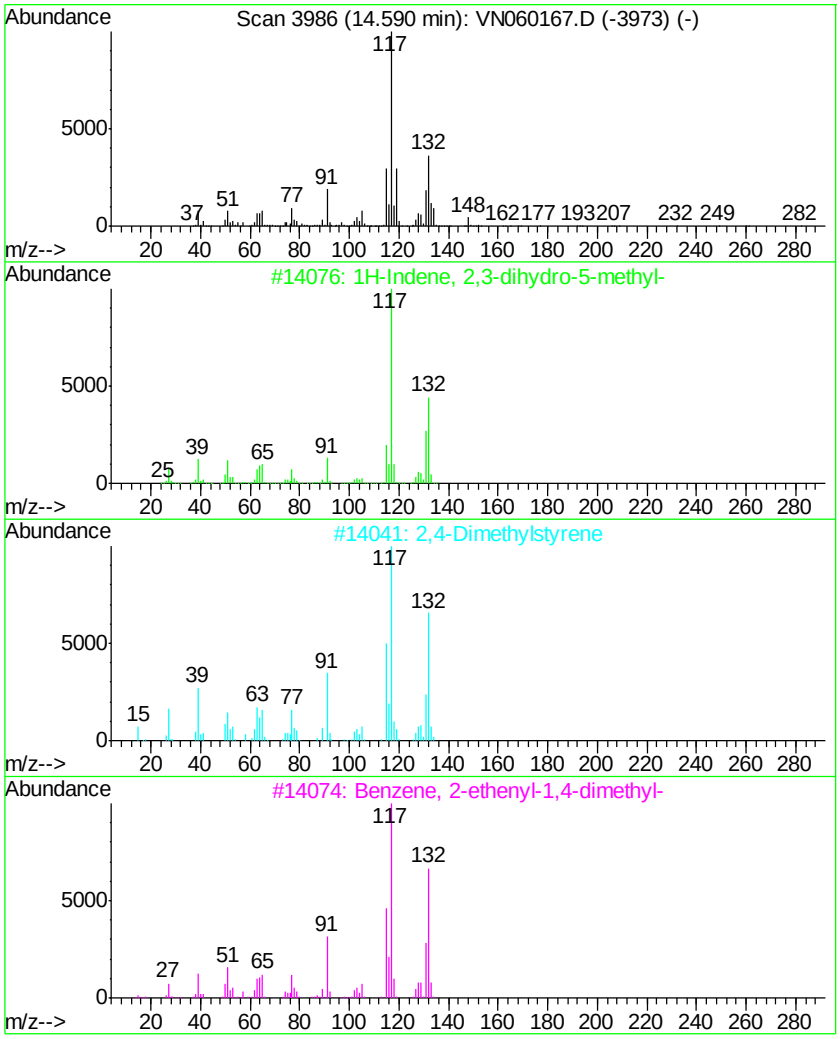
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

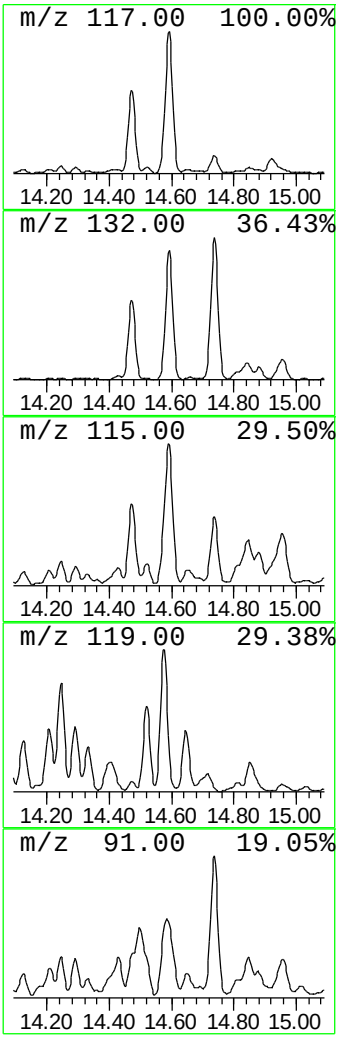
****

Peak Number 4 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.59	226.89 ug/l	4593120	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
2	2,4-Dimethylstyrene	132	C10H12	002234-20-0	90
3	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	90
4	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	89
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	86





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Operator : JC/MD
Sample : L1589-01 10X
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ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : SB3-9.5-10.0

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N021220W.M
Quant Title : SW846 8260

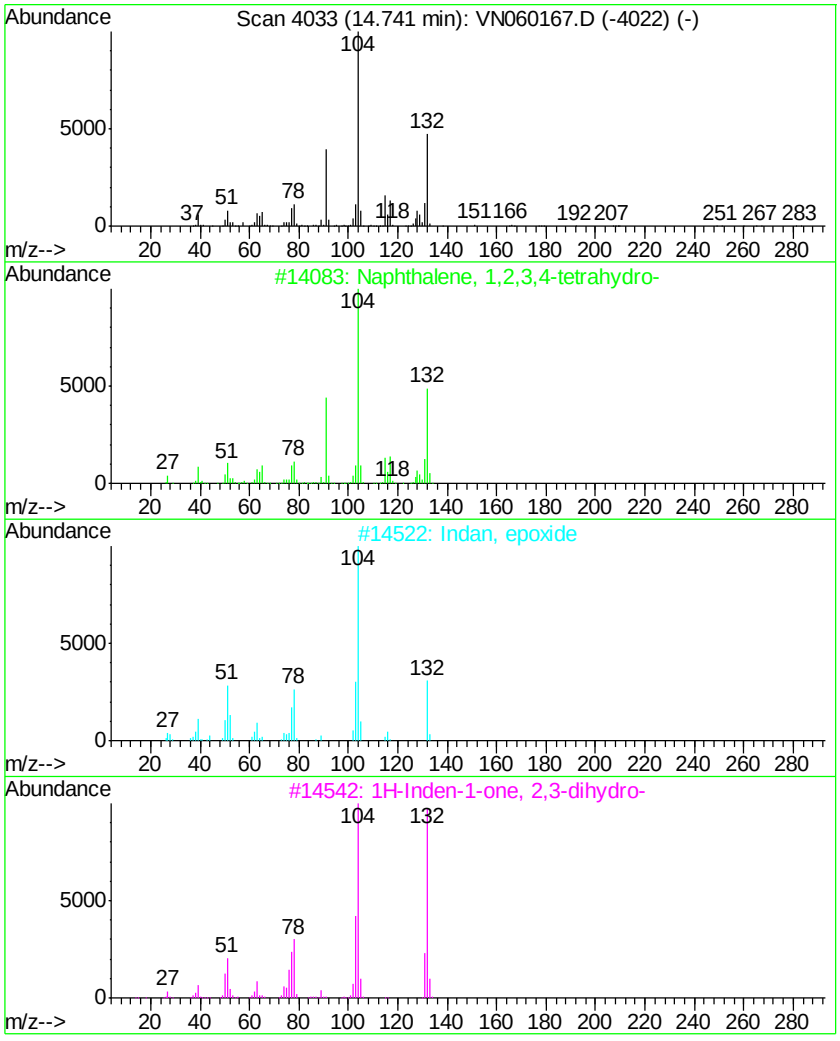
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

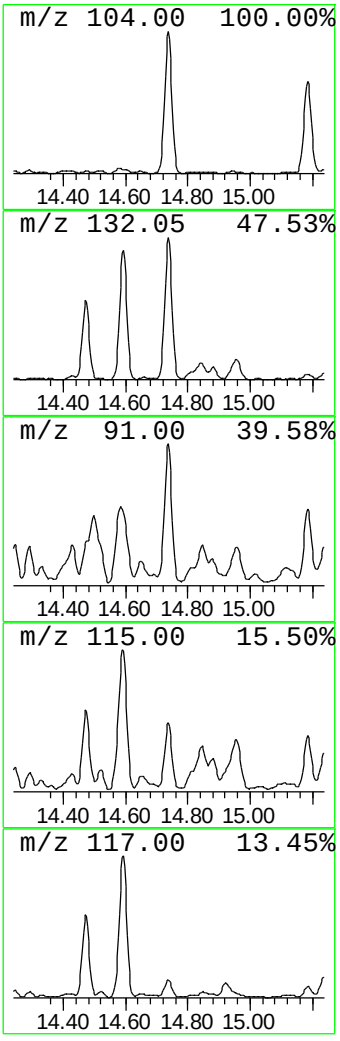
****

Peak Number 5 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.74	123.10 ug/l	2492000	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2	Indan, epoxide	132	C9H8O	000768-22-9	58
3	1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	42
4	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-92-1	42
5	Benzene, 3-cyclohexen-1-yl-	158	C12H14	004994-16-5	40





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN022020\
Data File : VN060167.D
Acq On : 20 Feb 2020 12:53
Operator : JC/MD
Sample : L1589-01 10X
Misc : 11.95g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : SB3-9.5-10.0

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N021220W.M
Quant Title : SW846 8260

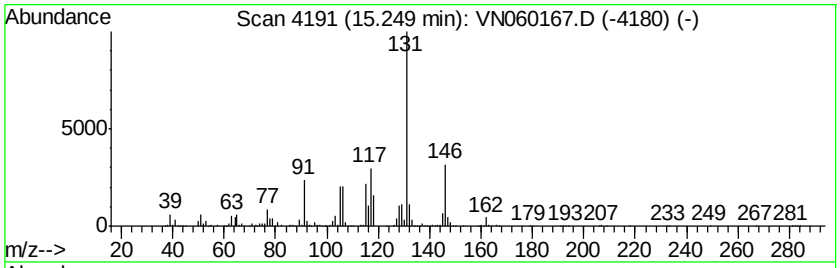
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

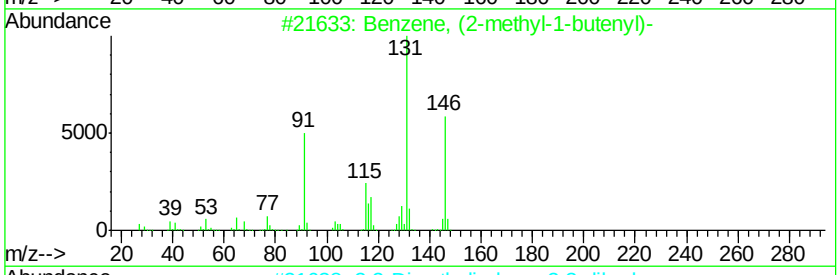
****

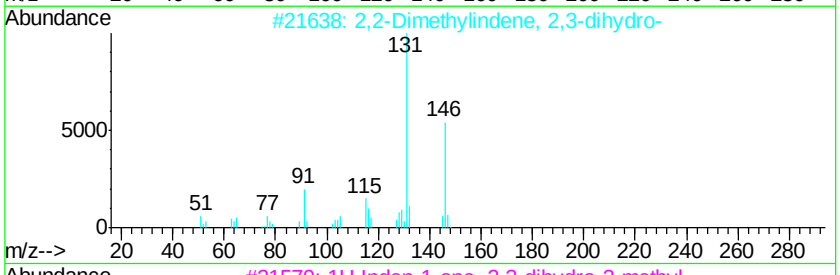
Peak Number 6 Benzene, (2-methyl-1-butenyl)- Concentration Rank 7

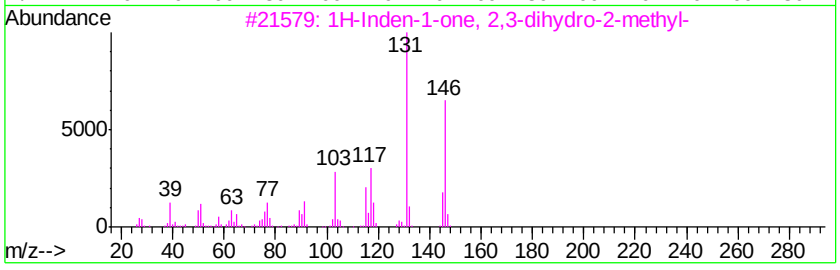
R.T.	EstConc	Area	Relative to ISTD	R.T.
15.25	133.00 ug/l	2692440	1,4-Dichlorobenzene-d4	13.34

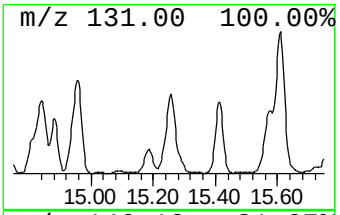
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	94
2	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	89
3	1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	81
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	81
5	Bicyclo[4.2.0]octa-1,3,5-triene,...	160	C12H16	078920-29-3	76

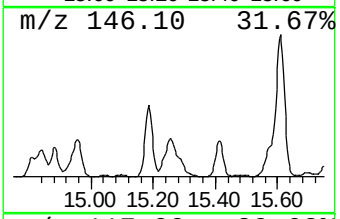


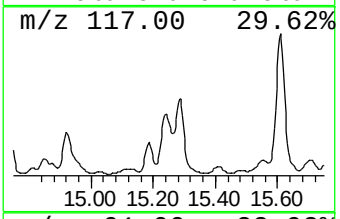


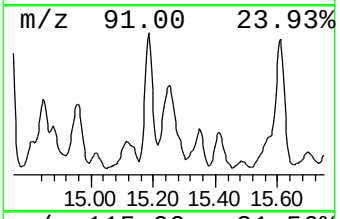


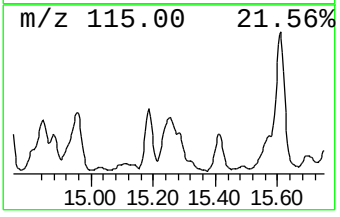












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Data File : VN060167.D
Acq On : 20 Feb 2020 12:53
Operator : JC/MD
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ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : SB3-9.5-10.0

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N021220W.M
Quant Title : SW846 8260

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

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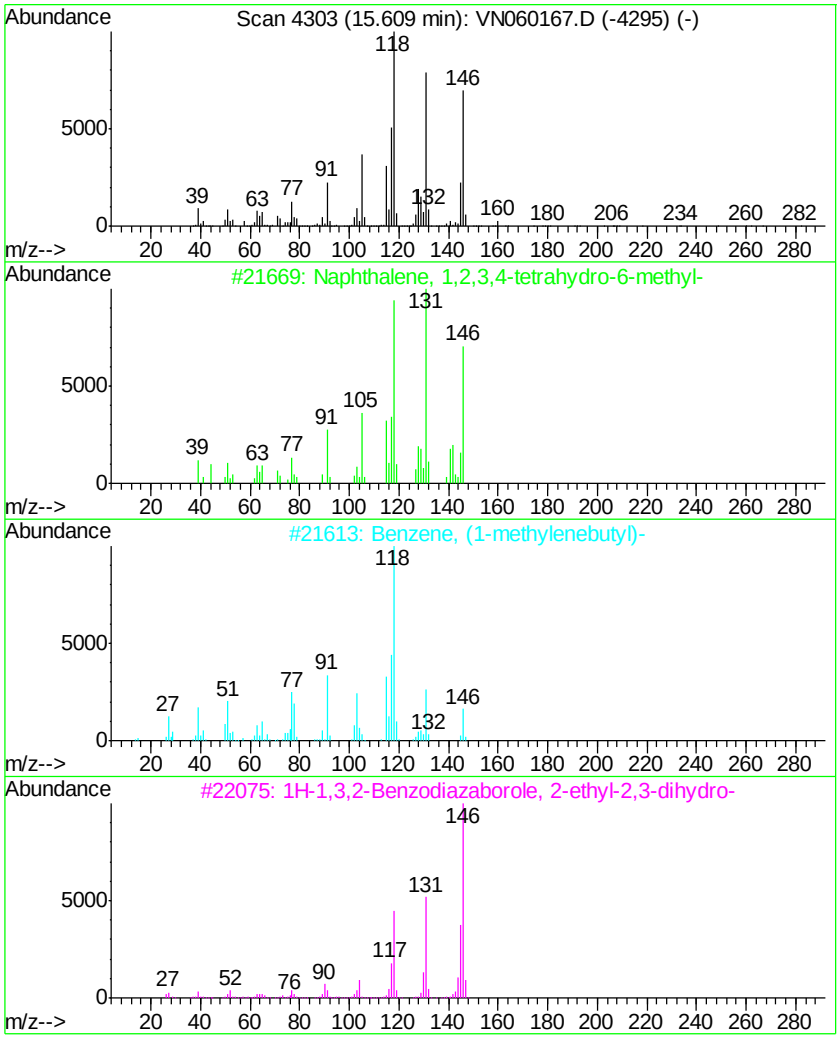
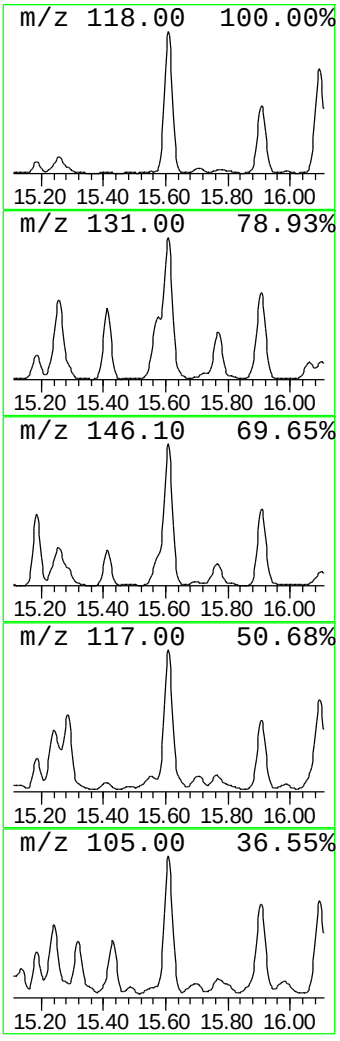
Peak Number 7 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.61	207.92 ug/l	4209180	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	91
2	Benzene, (1-methylenebutyl)-	146	C11H14	005676-32-4	72
3	1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	70
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	70
5	2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14	1000189-31-0	64





Library Search Compound Report

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Data File : VN060167.D
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Operator : JC/MD
Sample : L1589-01 10X
Misc : 11.95g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : SB3-9.5-10.0

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N021220W.M
Quant Title : SW846 8260

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

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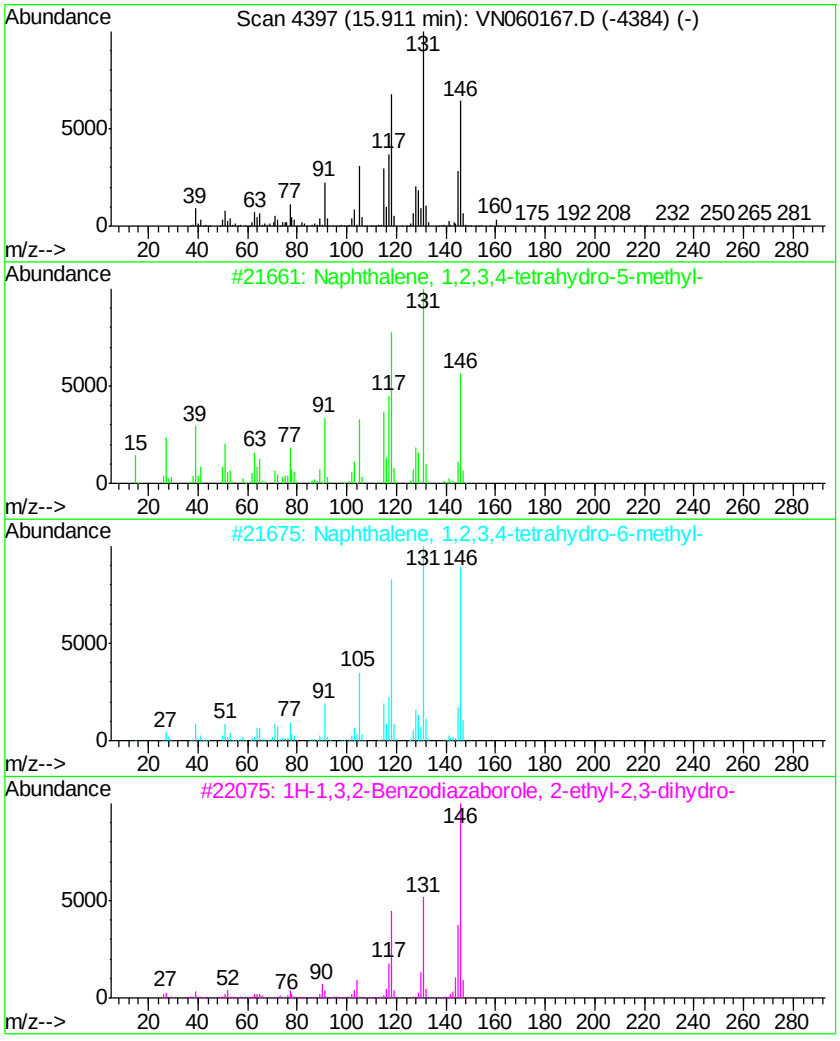
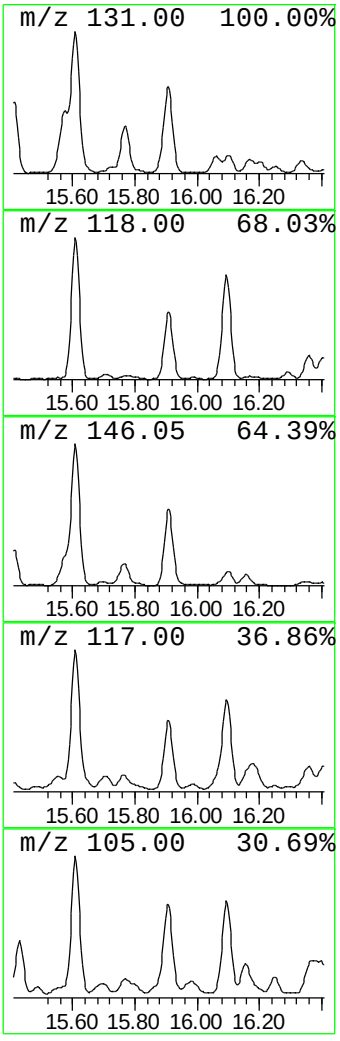
Peak Number 8 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.91	130.47 ug/l	2641210	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	94
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	87
3	1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	87
4	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	70
5	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	70





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ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : SB3-9.5-10.0

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N021220W.M
Quant Title : SW846 8260

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

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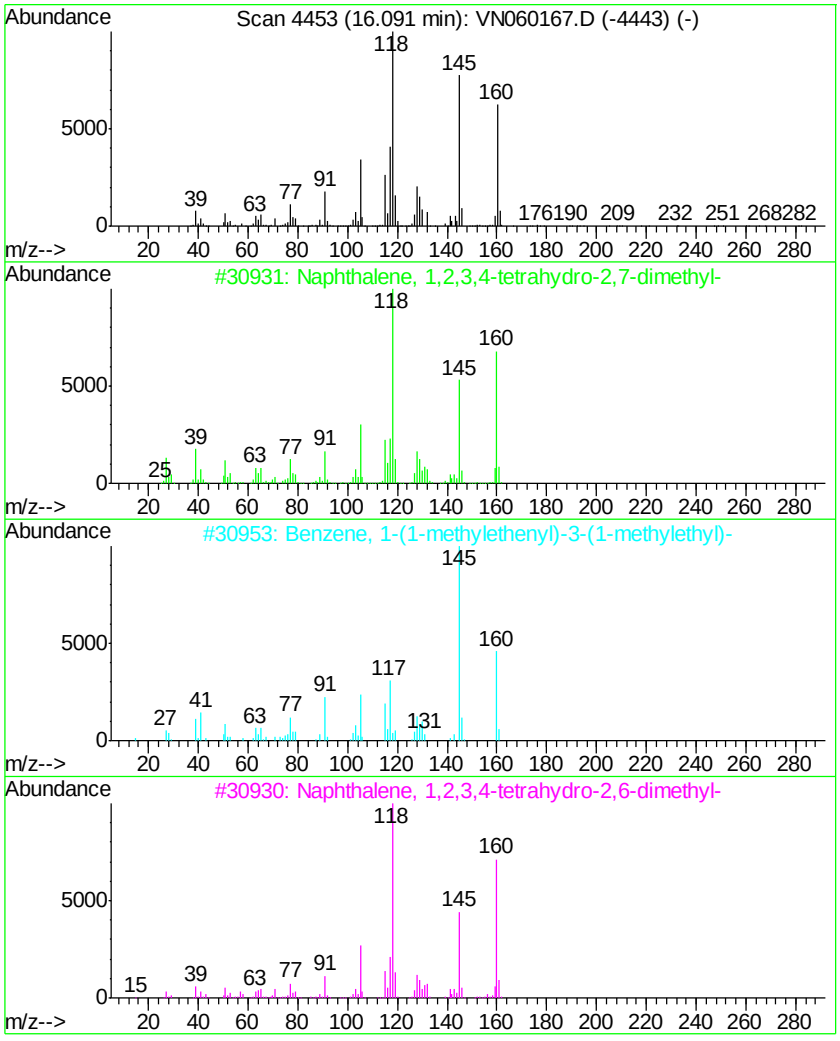
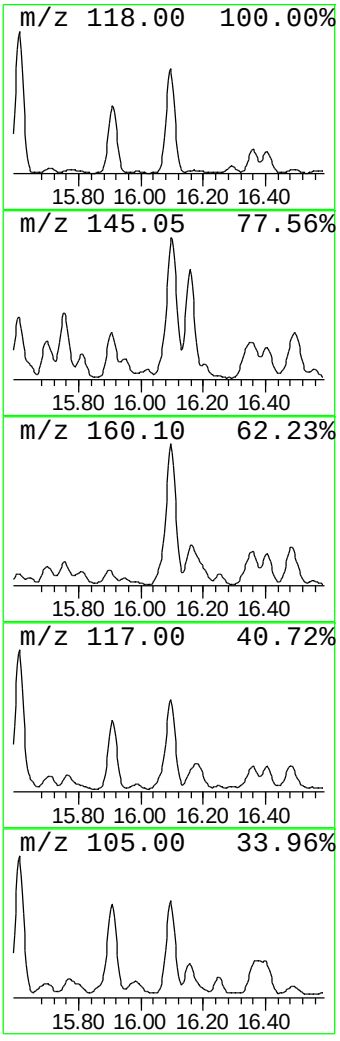
Peak Number 9 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.09	128.01 ug/l	2591450	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	89
2	Benzene, 1-(1-methylethenyl)-3-(...	160	C12H16	001129-29-9	70
3	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	007524-63-2	64
4	Benzene, 1,2,4-trimethyl-5-(1-me...	160	C12H16	054340-84-0	64
5	1H-Indene, 2,3-dihydro-1,1,3-tri...	160	C12H16	002613-76-5	60





Library Search Compound Report

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Quant Title : SW846 8260

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

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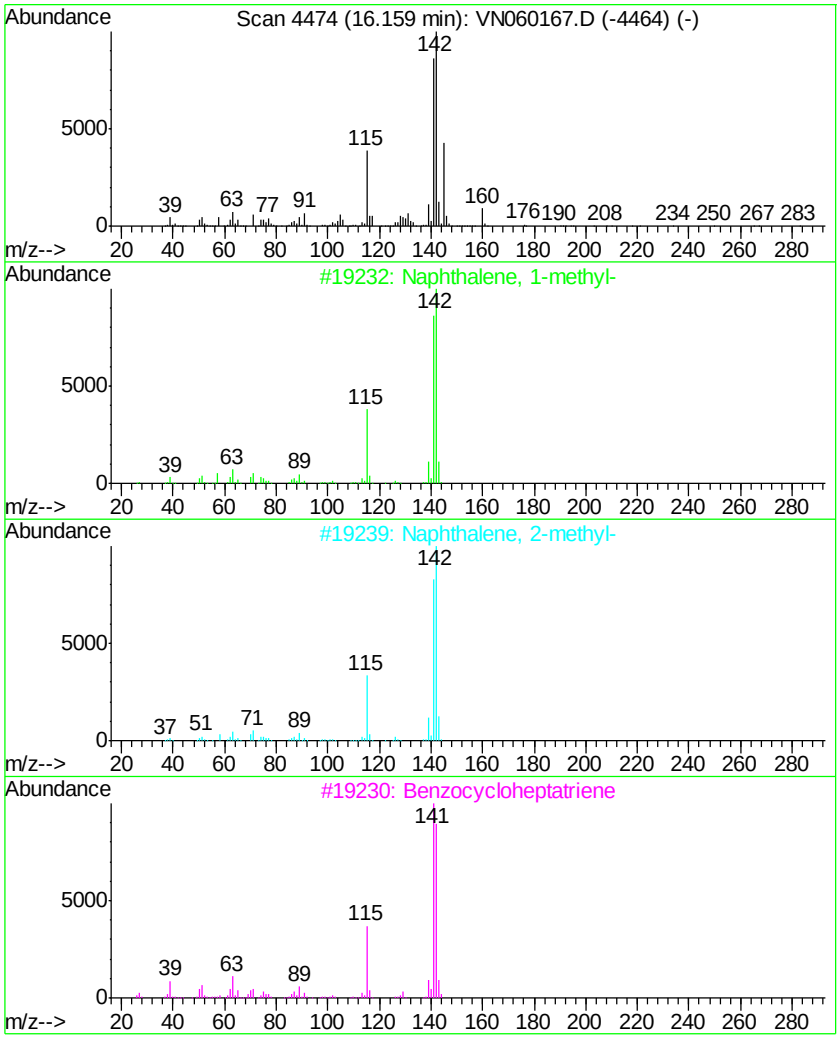
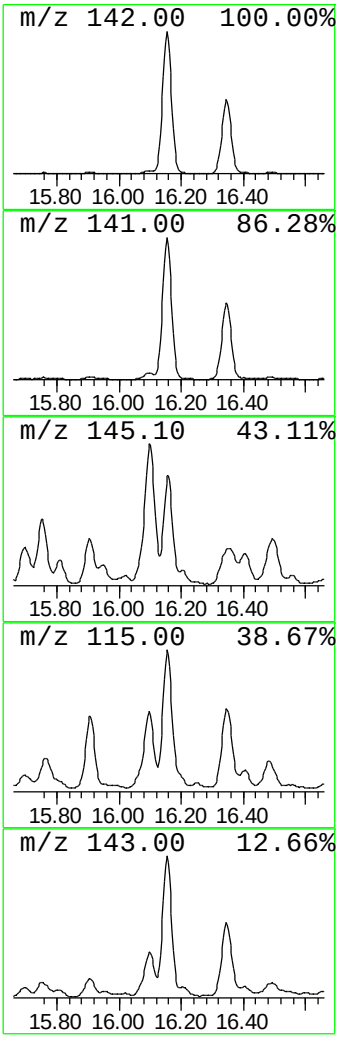
Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.16	141.25 ug/l	2859370	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3	Benzocycloheptatriene	142	C11H10	000264-09-5	76
4	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	76
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	49





Tentatively Identified Compound (LSC) summary

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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N021220W.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
Benzene, 1-ethyl-...	12.65	190.3	ug/l	3851560	4	13.34	1012200	50.0
Benzene, 1-methyl...	13.54	180.5	ug/l	3654690	4	13.34	1012200	50.0
unknown13.99	13.99	137.7	ug/l	2787530	4	13.34	1012200	50.0
1H-Indene, 2,3-di...	14.59	226.9	ug/l	4593120	4	13.34	1012200	50.0
Naphthalene, 1,2,...	14.74	123.1	ug/l	2492000	4	13.34	1012200	50.0
Benzene, (2-methy...	15.25	133.0	ug/l	2692440	4	13.34	1012200	50.0
Naphthalene, 1,2,...	15.61	207.9	ug/l	4209180	4	13.34	1012200	50.0
Naphthalene, 1,2,...	15.91	130.5	ug/l	2641210	4	13.34	1012200	50.0
Naphthalene, 1,2,...	16.09	128.0	ug/l	2591450	4	13.34	1012200	50.0
Naphthalene, 1-me...	16.16	141.3	ug/l	2859370	4	13.34	1012200	50.0