LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV070620\
Data File : VV017351.D
Acq On : 06 Jul 2020 14:53
Operator : SY/MD
Sample : L3154-18
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : H4298

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.1 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_V\METHOD\SOMVTR061720WMA.M
Title : TRACE VOA SOM01.0

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.116	3	8	18	rVB2	104156	98997	17.84%	1.376%
2	1.222	36	41	47	rVB	19707	16114	2.90%	0.224%
3	1.315	61	70	81	rBV3	166243	193845	34.93%	2.695%
4	1.364	81	85	90	rVV	10147	9444	1.70%	0.131%
5	1.409	90	99	100	rVV2	13110	18040	3.25%	0.251%
6	1.425	101	104	114	rVB2	18843	20429	3.68%	0.284%
7	1.579	136	152	161	rBV2	62180	92016	16.58%	1.279%
8	1.643	166	172	179	rBV2	19760	24438	4.40%	0.340%
9	1.775	209	213	218	rBV5	6067	6523	1.18%	0.091%
10	1.817	220	226	236	rVB	25093	33487	6.03%	0.466%
11	1.913	252	256	262	rVB2	6352	5991	1.08%	0.083%
12	1.994	276	281	288	rVB10	6949	8415	1.52%	0.117%
13	2.122	309	321	334	rBV	111837	168876	30.43%	2.348%
14	2.775	516	524	537	rVB4	19457	34894	6.29%	0.485%
15	2.978	577	587	596	rBV10	6614	15306	2.76%	0.213%
16	3.052	607	610	620	rVB10	6073	9276	1.67%	0.129%
17	3.534	749	760	772	rBV8	7160	15725	2.83%	0.219%
18	3.942	867	887	927	rBV2	170972	536078	96.61%	7.453%
19	4.376	1007	1022	1043	rBV	99531	246895	44.49%	3.433%
20	5.074	1221	1239	1266	rBV2	212873	554907	100.00%	7.715%
21	5.357	1319	1327	1331	rBV6	6799	11011	1.98%	0.153%
22	5.643	1402	1416	1440	rBV	207881	416316	75.02%	5.788%
23	5.971	1495	1518	1532	rBV	21138	69898	12.60%	0.972%
24	6.096	1542	1557	1580	rBV2	114967	244806	44.12%	3.404%
25	7.010	1829	1841	1860	rBV	57041	105892	19.08%	1.472%
26	7.338	1930	1943	1959	rBV	246675	424317	76.47%	5.900%
27	7.412	1959	1966	1976	rVB	13728	22418	4.04%	0.312%
28	7.646	2025	2039	2056	rBV	33392	65415	11.79%	0.910%
29	8.116	2175	2185	2221	rBV	237453	468286	84.39%	6.511%
30	8.874	2407	2421	2437	rBV	311247	509349	91.79%	7.082%
31	9.035	2461	2471	2481	rVB	28095	45329	8.17%	0.630%
32	9.164	2495	2511	2527	rVB	17266	35297	6.36%	0.491%
33	9.566	2632	2636	2646	rVB9	4417	6417	1.16%	0.089%
34	9.955	2747	2757	2762	rBV8	5196	7456	1.34%	0.104%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV070620\
Data File : VV017351.D
Acq On : 06 Jul 2020 14:53
Operator : SY/MD
Sample : L3154-18
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : H4298

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.1 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_V\METHOD\SOMVTR061720WMA.M
Title : TRACE VOA SOM01.0

35	10.238	2833	2845	2862	rBV	95014	153372	27.64%	2.132%
36	10.710	2987	2992	3001	rVB	4466	6227	1.12%	0.087%
37	11.270	3155	3166	3183	rBV	304611	482851	87.01%	6.713%
38	11.556	3243	3255	3258	rBV8	8433	11951	2.15%	0.166%
39	11.646	3271	3283	3300	rBV	273883	442720	79.78%	6.155%
40	11.768	3317	3321	3332	rVB10	4264	6169	1.11%	0.086%
41	11.842	3334	3344	3352	rBV	16166	24666	4.45%	0.343%
42	12.138	3428	3436	3445	rBV3	11257	16979	3.06%	0.236%
43	12.434	3520	3528	3538	rVB	17021	25792	4.65%	0.359%
44	12.489	3539	3545	3555	rVB6	5846	8814	1.59%	0.123%
45	12.704	3603	3612	3615	rBV6	5342	8078	1.46%	0.112%
46	12.903	3664	3674	3684	rVB3	16237	25780	4.65%	0.358%
47	12.977	3690	3697	3705	rVB3	5408	7920	1.43%	0.110%
48	13.064	3719	3724	3732	rBV3	3663	5712	1.03%	0.079%
49	13.292	3789	3795	3801	rBV7	6646	8716	1.57%	0.121%
50	13.331	3801	3807	3814	rVB7	5466	7925	1.43%	0.110%
51	13.636	3891	3902	3914	rBV	101065	158951	28.64%	2.210%
52	13.726	3916	3930	3947	rVV	116825	196362	35.39%	2.730%
53	13.797	3948	3952	3961	rVB10	4558	6300	1.14%	0.088%
54	14.109	4040	4049	4052	rBV3	17787	24138	4.35%	0.336%
55	14.138	4053	4058	4072	rVB3	27092	47635	8.58%	0.662%
56	14.247	4081	4092	4103	rBV5	40082	71509	12.89%	0.994%
57	14.318	4103	4114	4127	rVV5	50117	101499	18.29%	1.411%
58	14.382	4127	4134	4147	rVB10	13371	22869	4.12%	0.318%
59	14.476	4151	4163	4169	rBV4	14091	27466	4.95%	0.382%
60	14.521	4169	4177	4182	rBV9	10075	13130	2.37%	0.183%
61	14.595	4194	4200	4202	rBV7	7236	6879	1.24%	0.096%
62	14.640	4206	4214	4217	rBV6	23905	32012	5.77%	0.445%
63	14.710	4227	4236	4249	rVB4	83995	158763	28.61%	2.207%
64	14.787	4252	4260	4270	rBV4	23556	39594	7.14%	0.550%
65	14.852	4270	4280	4288	rVB5	13033	24207	4.36%	0.337%
66	14.900	4290	4295	4296	rBV4	9371	7429	1.34%	0.103%
67	14.935	4303	4306	4322	rVB10	20330	36066	6.50%	0.501%
68	15.125	4359	4365	4372	rVB4	16086	22699	4.09%	0.316%
69	15.177	4372	4381	4396	rVB4	25316	56916	10.26%	0.791%
70	15.270	4406	4410	4417	rVB8	5668	6345	1.14%	0.088%
71	15.340	4417	4432	4437	rBV4	31756	63045	11.36%	0.877%

LSC Area Percent Report

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Acq On : 06 Jul 2020 14:53
Operator : SY/MD
Sample : L3154-18
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : H4298

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.1 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_V\METHOD\SOMVTR061720WMA.M
Title : TRACE VOA SOM01.0

72	15.450	4455	4466	4479	rBV8	42575	80691	14.54%	1.122%
73	15.524	4482	4489	4500	rVB3	17087	26926	4.85%	0.374%
74	15.585	4500	4508	4513	rBV5	15621	25498	4.60%	0.355%
75	15.701	4537	4544	4551	rBV8	16853	31294	5.64%	0.435%
76	15.746	4553	4558	4563	rVB4	13911	18890	3.40%	0.263%
77	15.836	4580	4586	4593	rBV8	13042	22687	4.09%	0.315%
78	16.086	4656	4664	4673	rVB8	12800	23388	4.21%	0.325%
79	16.138	4674	4680	4687	rVB10	7197	10211	1.84%	0.142%
80	16.176	4687	4692	4696	rBV7	6442	7153	1.29%	0.099%
81	16.209	4699	4702	4711	rVV8	7867	12104	2.18%	0.168%
82	16.260	4711	4718	4728	rVB8	10763	17511	3.16%	0.243%
83	16.498	4787	4792	4802	rVB8	4023	6641	1.20%	0.092%

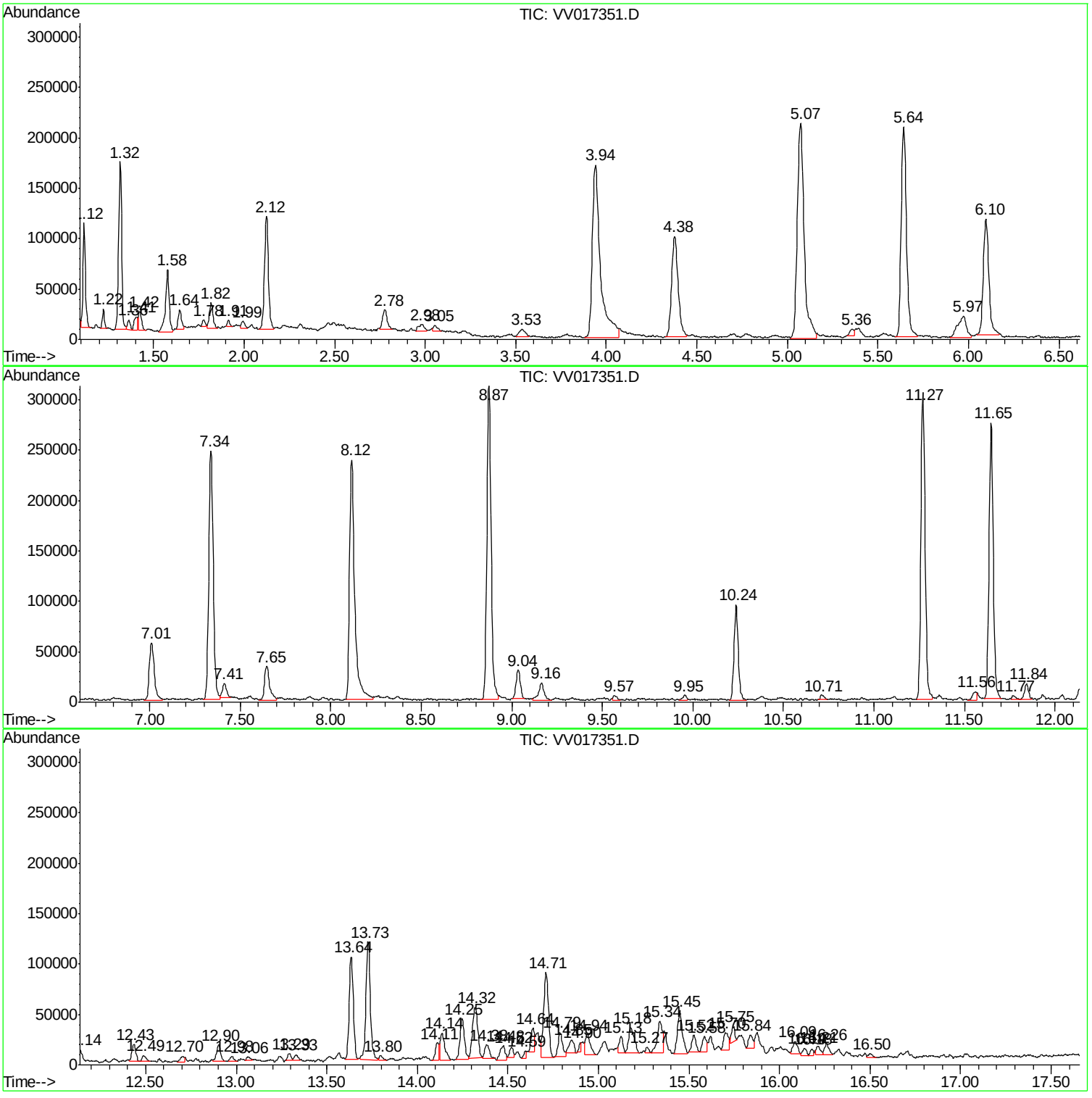
Sum of corrected areas: 7192383

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV070620\
Data File : VV017351.D
Acq On : 06 Jul 2020 14:53
Operator : SY/MD
Sample : L3154-18
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : H4298

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR061720WMA.M
Quant Title : TRACE VOA SOM01.0

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



Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV070620\
Data File : VV017351.D
Acq On : 06 Jul 2020 14:53
Operator : SY/MD
Sample : L3154-18
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : H4298

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR061720WMA.M
Quant Title : TRACE VOA SOM01.0

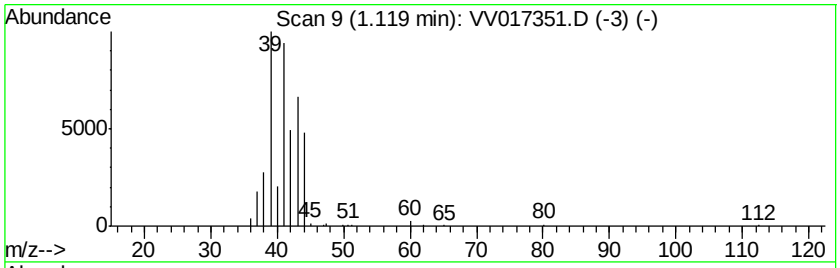
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

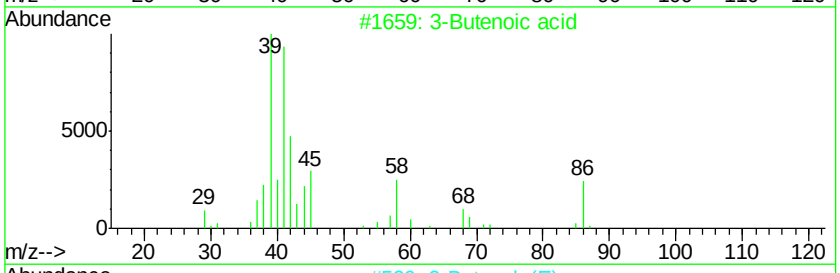
****

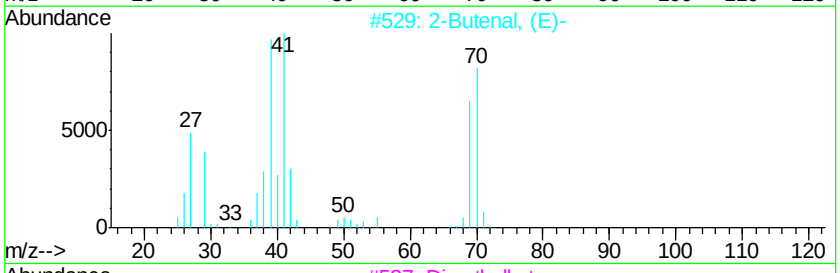
Peak Number 1 3-Butenoic acid Concentration Rank 5

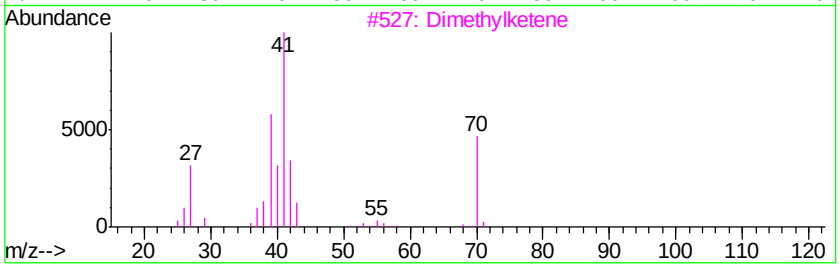
R.T.	EstConc	Area	Relative to ISTD	R.T.
1.12	1.19 ug/L	98997	1,4-Difluorobenzene	5.64

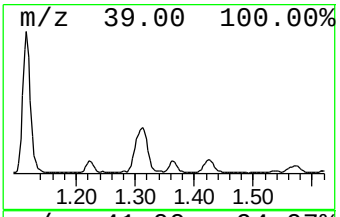
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Butenoic acid	86	C4H6O2	000625-38-7	64
2	2-Butenal, (E)-	70	C4H6O	000123-73-9	56
3	Dimethylketene	70	C4H6O	000598-26-5	40
4	1-Propyne, 3-chloro-	74	C3H3Cl	000624-65-7	39
5	Thiocyanic acid, 2-propynyl ester	97	C4H3NS	024309-48-6	39

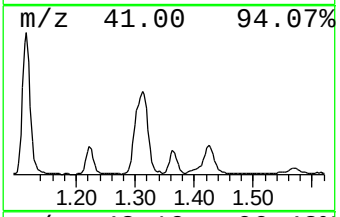


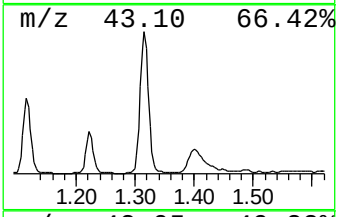


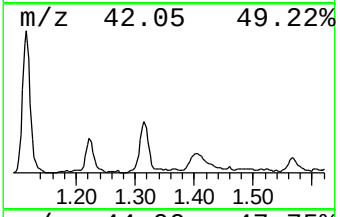


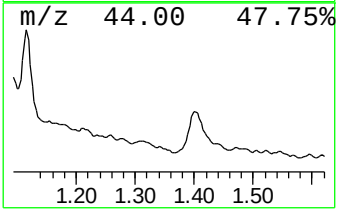












Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV070620\
Data File : VV017351.D
Acq On : 06 Jul 2020 14:53
Operator : SY/MD
Sample : L3154-18
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : H4298

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR061720WMA.M
Quant Title : TRACE VOA SOM01.0

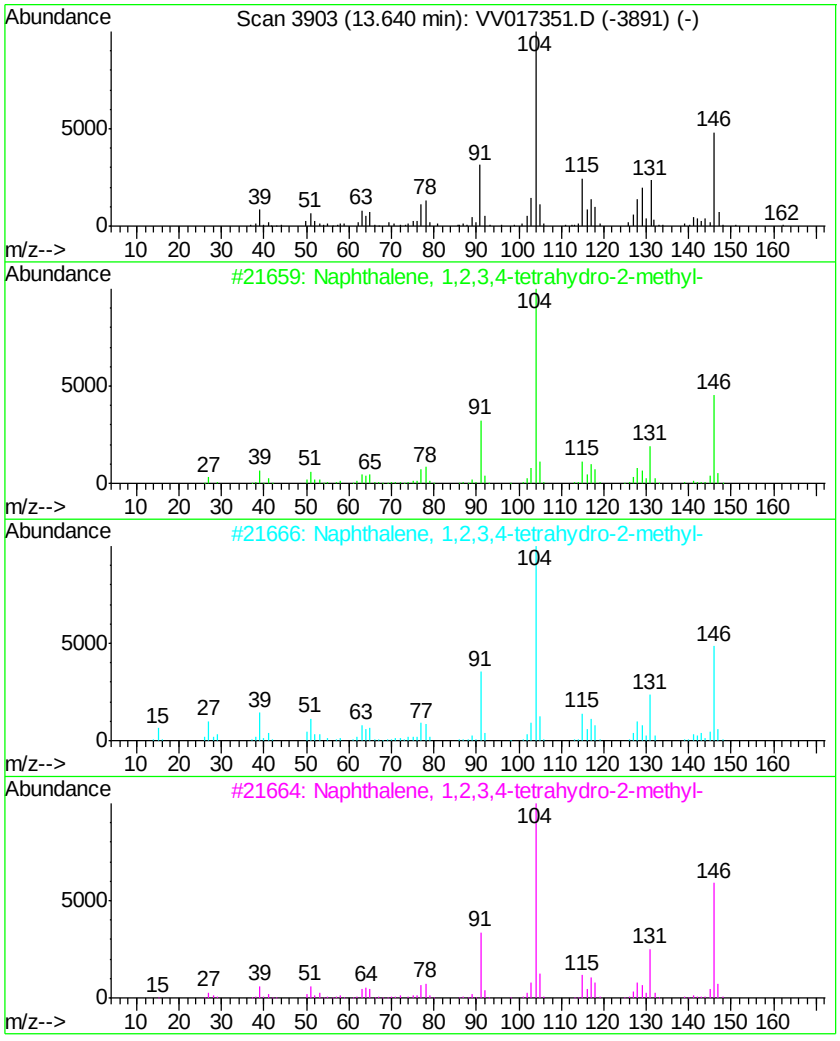
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

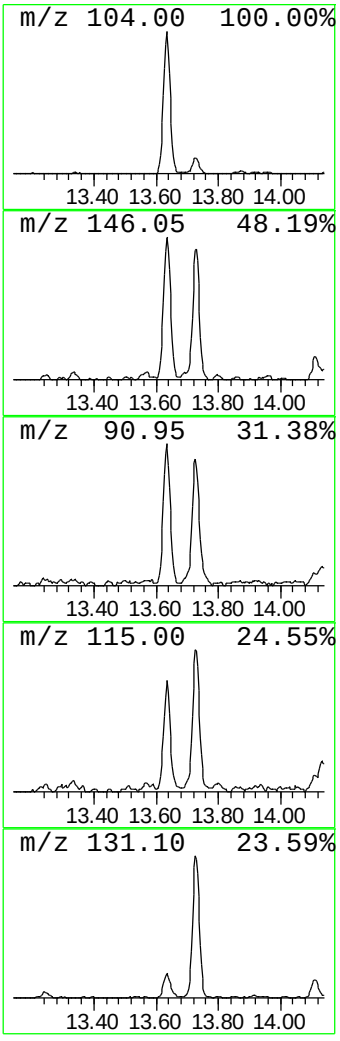
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.64	1.65 ug/L	158951	1,4-Dichlorobenzene-d4	11.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	91
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	91
3	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	74
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	74
5	Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV070620\
Data File : VV017351.D
Acq On : 06 Jul 2020 14:53
Operator : SY/MD
Sample : L3154-18
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : H4298

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR061720WMA.M
Quant Title : TRACE VOA SOM01.0

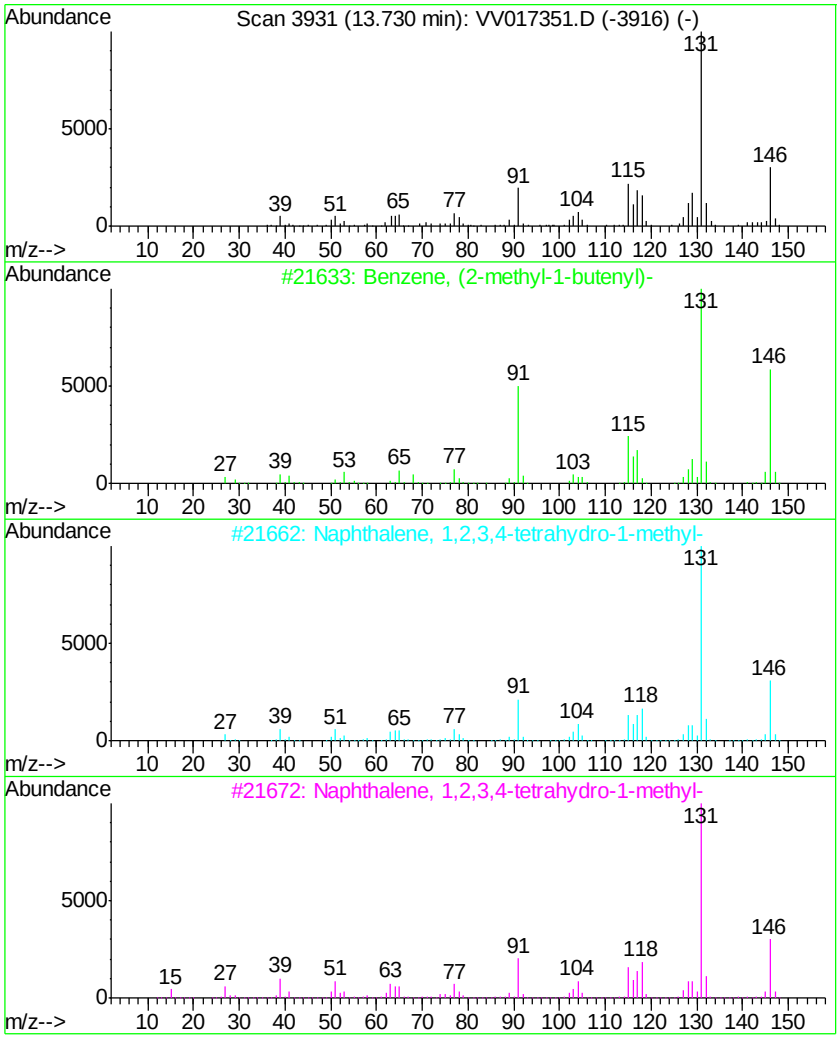
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

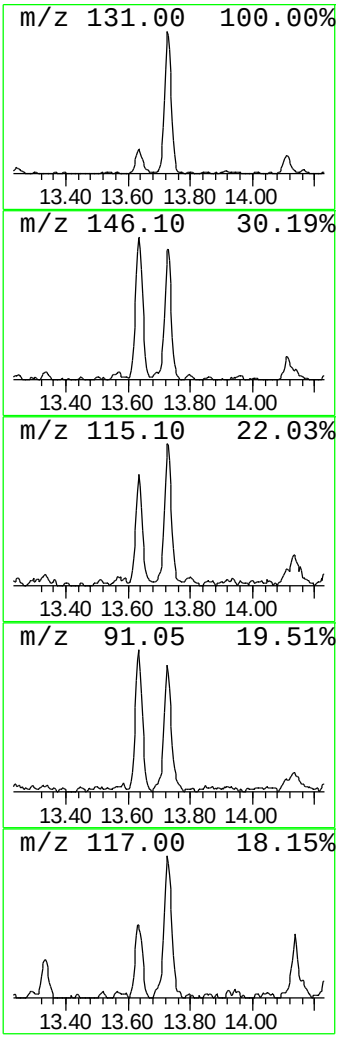
****

Peak Number 4 Benzene, (2-methyl-1-butenyl)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.73	2.03 ug/L	196362	1,4-Dichlorobenzene-d4	11.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	94
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	93
3	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	93
4	Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	90
5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	90





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV070620\
Data File : VV017351.D
Acq On : 06 Jul 2020 14:53
Operator : SY/MD
Sample : L3154-18
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : H4298

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR061720WMA.M
Quant Title : TRACE VOA SOM01.0

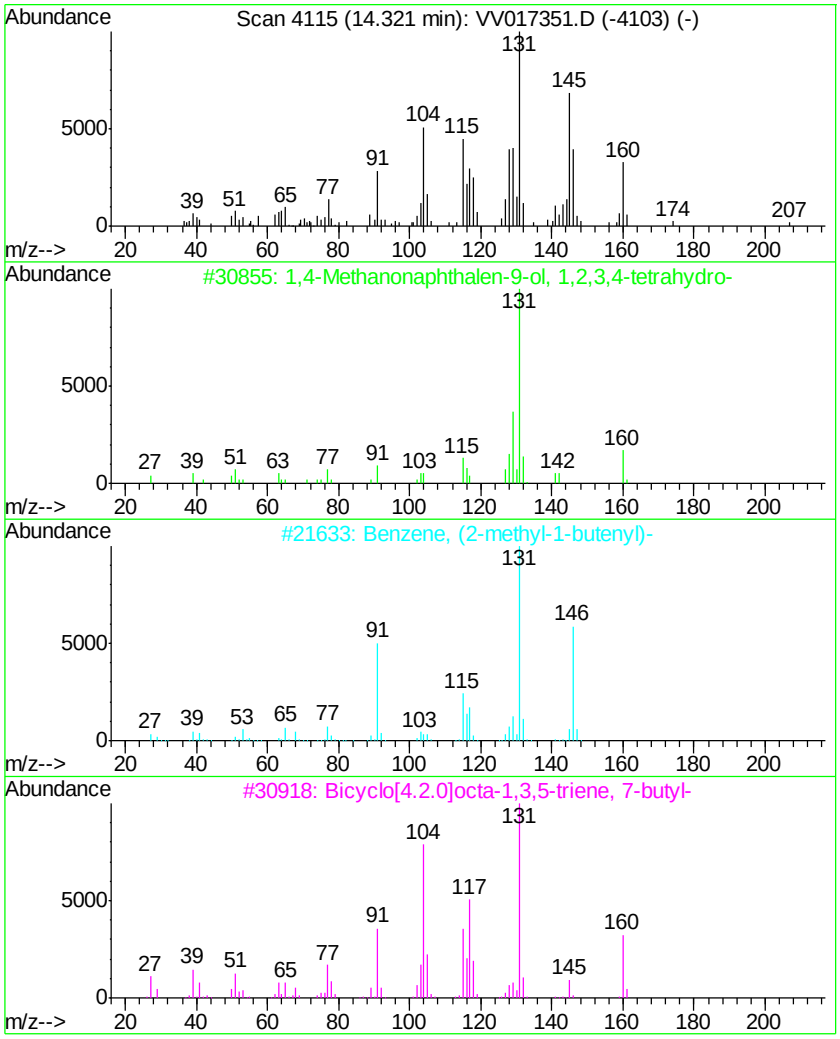
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

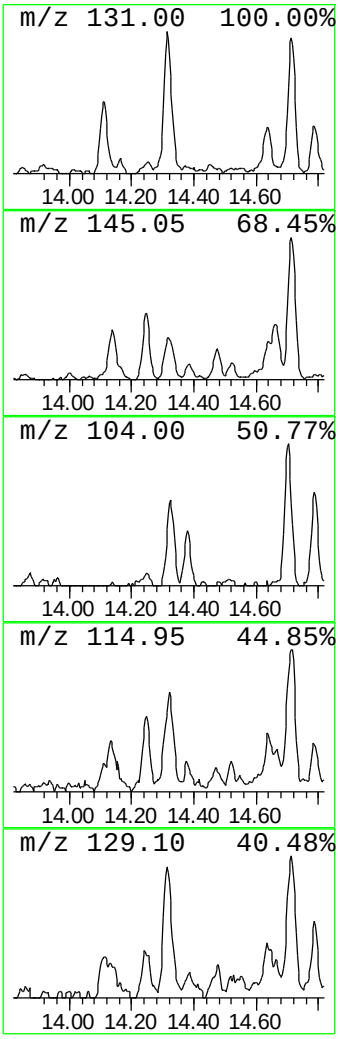
****

Peak Number 6 1,4-Methanonaphthalen-9-ol,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.32	1.05 ug/L	101499	1,4-Dichlorobenzene-d4	11.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Methanonaphthalen-9-ol, 1,2,...	160	C11H12O	055255-94-2	64
2	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	55
3	Bicyclo[4.2.0]octa-1,3,5-triene,...	160	C12H16	078920-29-3	55
4	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	55
5	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	49





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV070620\
Data File : VV017351.D
Acq On : 06 Jul 2020 14:53
Operator : SY/MD
Sample : L3154-18
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : H4298

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR061720WMA.M
Quant Title : TRACE VOA SOM01.0

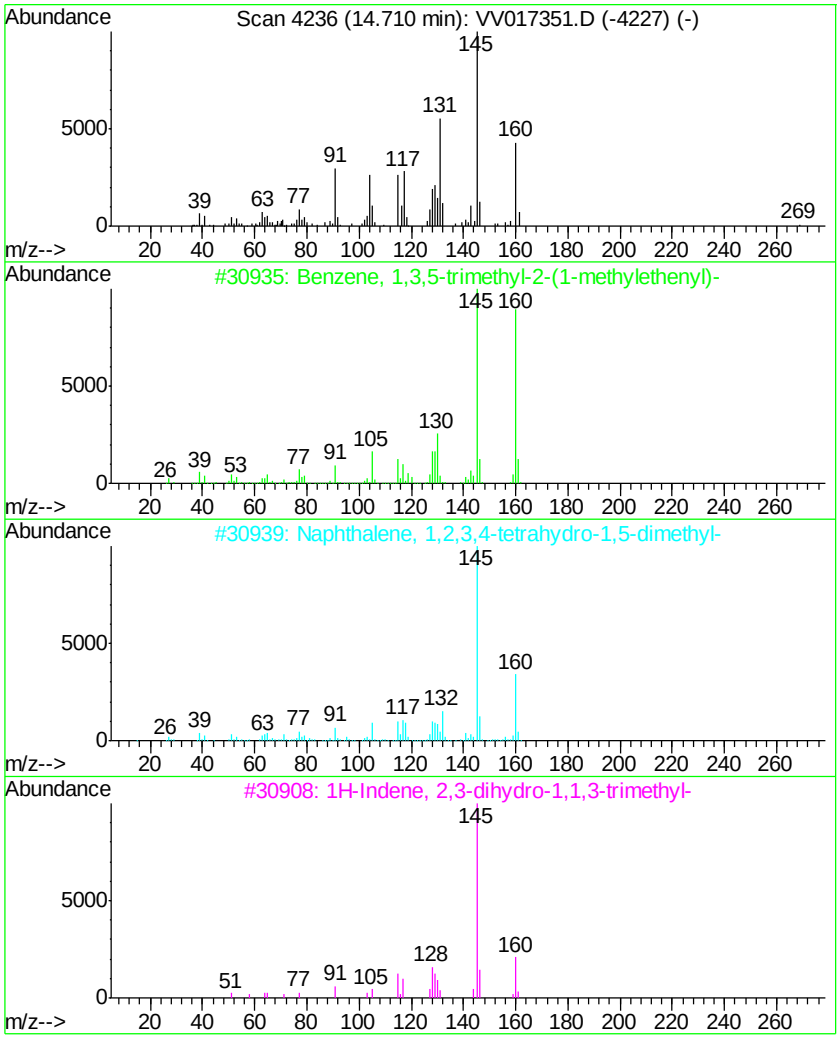
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

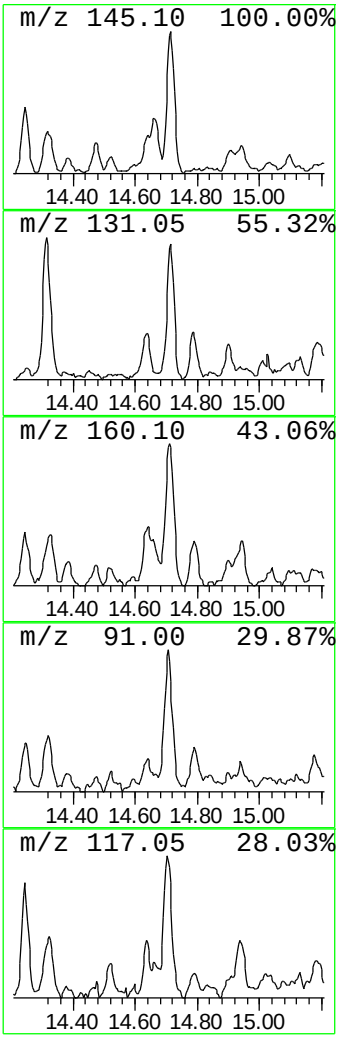
****

Peak Number 7 Benzene, 1,3,5-trimethyl-2-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.71	1.64 ug/L	158763	1,4-Dichlorobenzene-d4	11.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	70
2	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-91-0	70
3	1H-Indene, 2,3-dihydro-1,1,3-tri...	160	C12H16	002613-76-5	70
4	Benzene, (2,2-dimethyl-1-methyle...	160	C12H16	005676-29-9	64
5	Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	64





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV070620\
Data File : VV017351.D
Acq On : 06 Jul 2020 14:53
Operator : SY/MD
Sample : L3154-18
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : H4298

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR061720WMA.M
Quant Title : TRACE VOA SOM01.0

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

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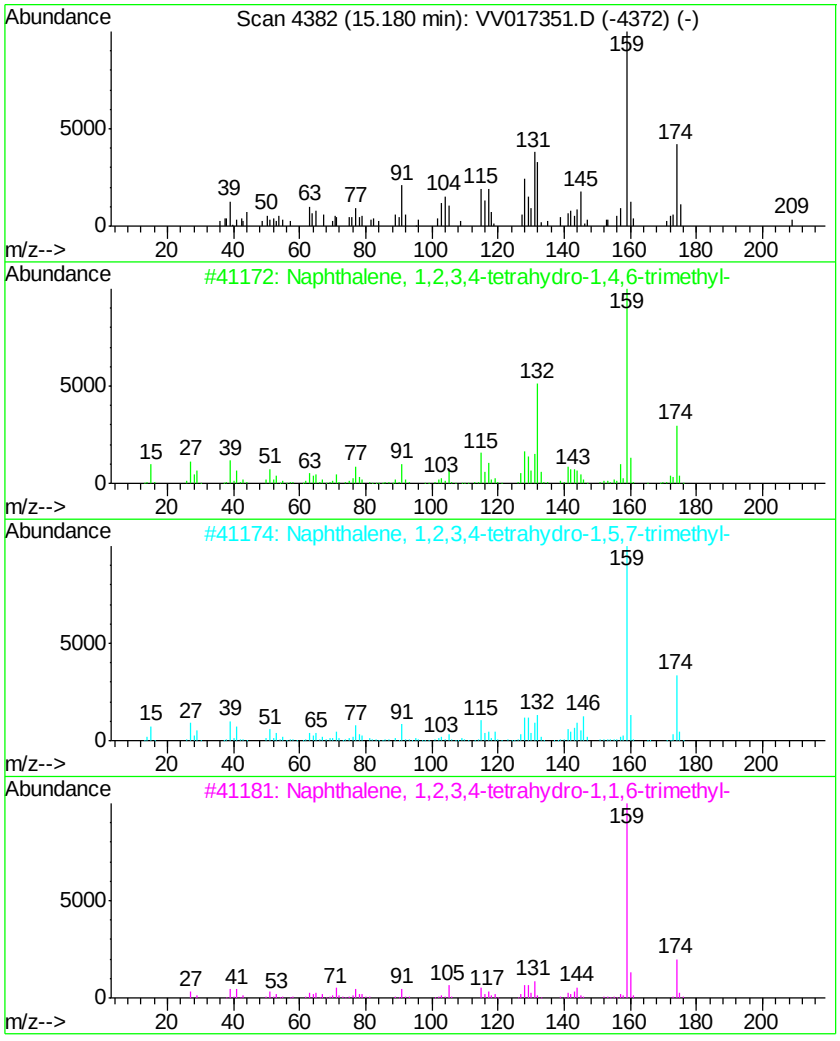
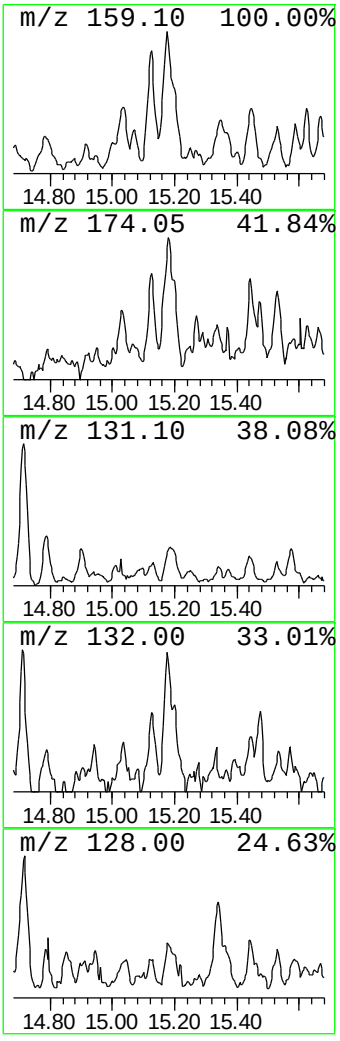
Peak Number 8 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.18	0.59 ug/L	56916	1,4-Dichlorobenzene-d4	11.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	022824-32-4	87
2	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	021693-55-0	74
3	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	000475-03-6	72
4	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	030316-36-0	70
5	1H-Inden-1-one, 2,3-dihydro-3,4,...	174	C12H14O	035322-84-0	70





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV070620\
Data File : VV017351.D
Acq On : 06 Jul 2020 14:53
Operator : SY/MD
Sample : L3154-18
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : H4298

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR061720WMA.M
Quant Title : TRACE VOA SOM01.0

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

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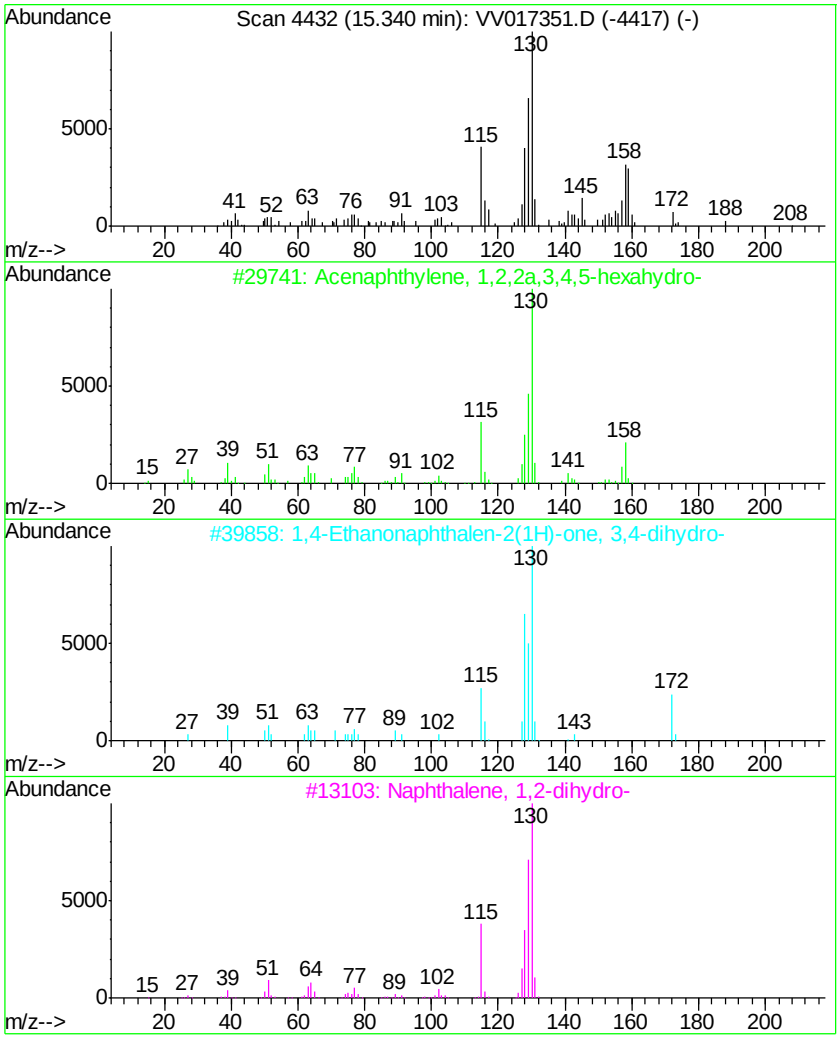
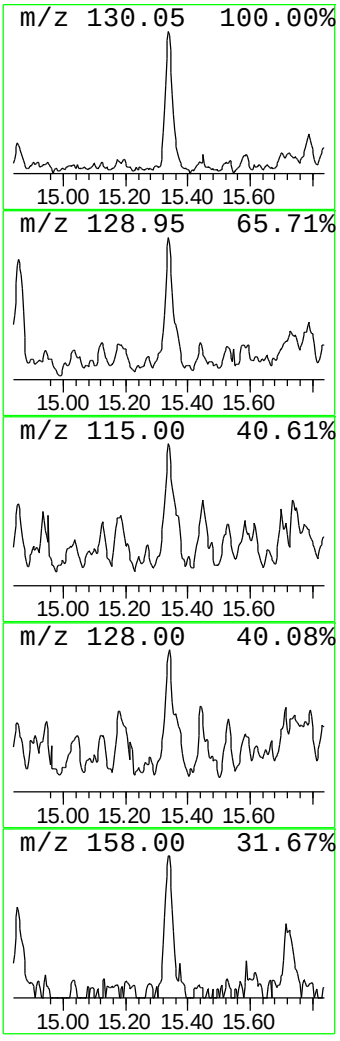
Peak Number 9 Acenaphthylene, 1,2,2a,3,4,... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.34	0.65 ug/L	63045	1,4-Dichlorobenzene-d4	11.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acenaphthylene, 1,2,2a,3,4,5-hex...	158	C12H14	000480-72-8	68
2	1,4-Ethanonaphthalen-2(1H)-one, ...	172	C12H12O	013153-76-9	60
3	Naphthalene, 1,2-dihydro-	130	C10H10	000447-53-0	60
4	Naphthalene, 1,2-dihydro-	130	C10H10	000447-53-0	60
5	1H-Indene-3-ethanamine	159	C11H13N	015940-39-3	58





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV070620\
Data File : VV017351.D
Acq On : 06 Jul 2020 14:53
Operator : SY/MD
Sample : L3154-18
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : H4298

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR061720WMA.M
Quant Title : TRACE VOA SOM01.0

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

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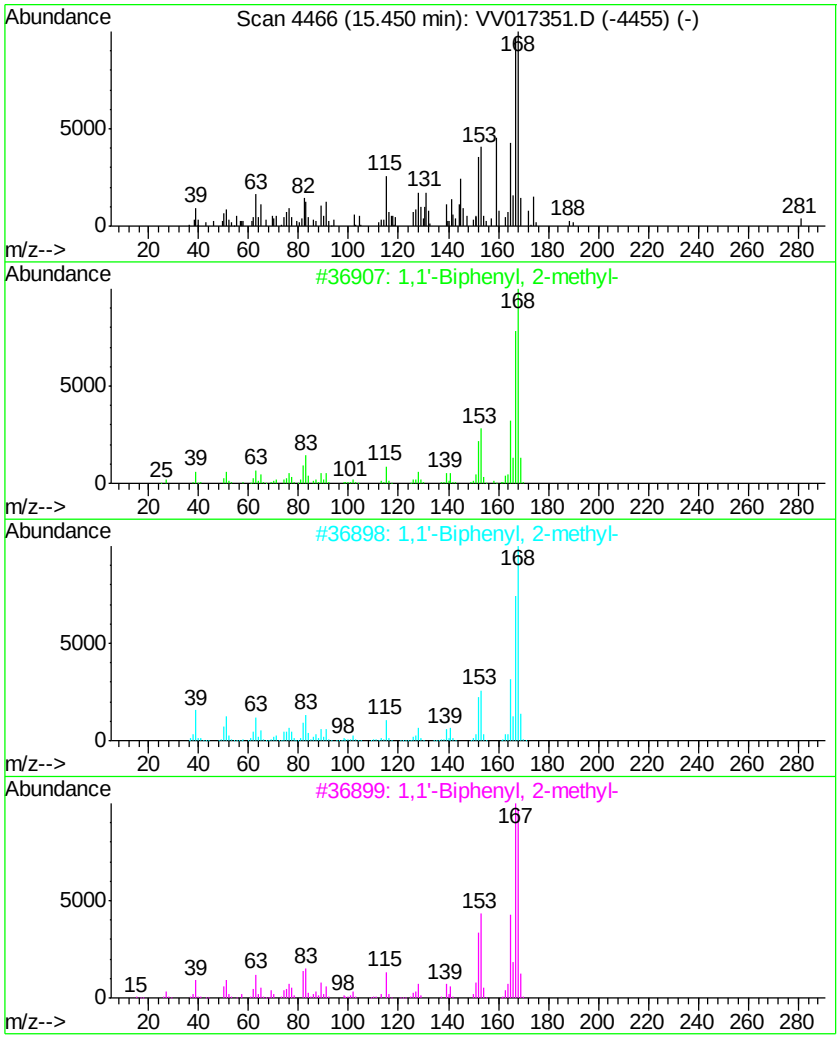
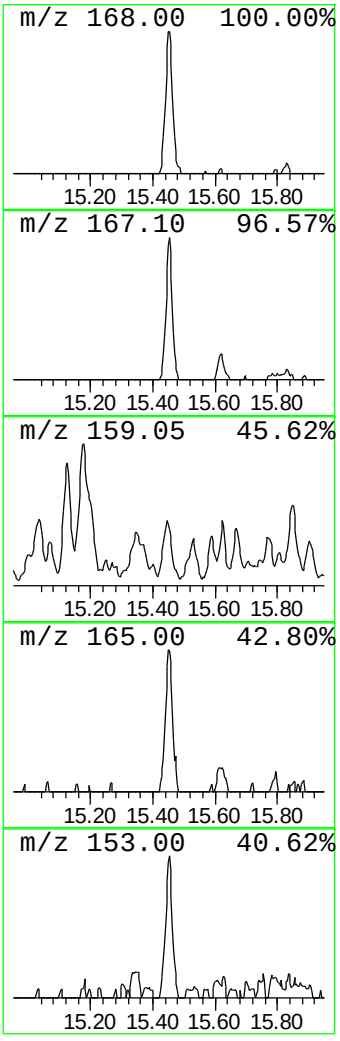
Peak Number 10 1,1'-Biphenyl, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.45	0.84 ug/L	80691	1,4-Dichlorobenzene-d4	11.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	96
2	1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	93
3	1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	93
4	1,1'-Biphenyl, 3-methyl-	168	C13H12	000643-93-6	83
5	Diphenylmethane	168	C13H12	000101-81-5	83





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV070620\
Data File : VV017351.D
Acq On : 06 Jul 2020 14:53
Operator : SY/MD
Sample : L3154-18
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : H4298

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR061720WMA.M
Quant Title : TRACE VOA SOM01.0

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
3-Butenoic acid	1.12	1.2	ug/L	98997	1	5.64	416316	5.0
Naphthalene, 1,2,...	13.64	1.6	ug/L	158951	3	11.27	482851	5.0
Benzene, (2-methy...	13.73	2.0	ug/L	196362	3	11.27	482851	5.0
1,4-Methanonaphth...	14.32	1.1	ug/L	101499	3	11.27	482851	5.0
Benzene, 1,3,5-tr...	14.71	1.6	ug/L	158763	3	11.27	482851	5.0
Naphthalene, 1,2,...	15.18	0.6	ug/L	56916	3	11.27	482851	5.0
Acenaphthylene, 1...	15.34	0.7	ug/L	63045	3	11.27	482851	5.0
1,1'-Biphenyl, 2-...	15.45	0.8	ug/L	80691	3	11.27	482851	5.0