LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122923\
Data File : VX039537.D
Acq On : 29 Dec 2023 12:04
Operator : JC/MD
Sample : 06017-01
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-7.5-8.5

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_X\Method\82X121323W.M
Title : SW846 8260

Signal : TIC: VX039537.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.391	694	706	715	rBV	58657	189140	19.75%	0.893%
2	5.489	715	722	724	rVV2	30377	76498	7.99%	0.361%
3	5.544	725	731	761	rVB	103603	368614	38.50%	1.739%
4	5.806	763	774	789	rBV2	37438	109060	11.39%	0.515%
5	5.958	789	799	807	rBV	75066	215682	22.53%	1.018%
6	6.031	807	811	820	rVV2	16719	41962	4.38%	0.198%
7	6.232	834	844	854	rVV2	91882	287056	29.98%	1.355%
8	6.598	893	904	916	rBV2	32990	84437	8.82%	0.398%
9	6.757	921	930	939	rBV	189400	469835	49.07%	2.217%
10	7.366	1019	1030	1040	rVV4	24683	69184	7.23%	0.326%
11	7.488	1042	1050	1055	rVV	39307	80452	8.40%	0.380%
12	7.555	1055	1061	1072	rVV	50130	120325	12.57%	0.568%
13	7.769	1088	1096	1104	rVB4	18603	38817	4.05%	0.183%
14	7.976	1120	1130	1141	rVB	107184	230133	24.03%	1.086%
15	8.098	1141	1150	1158	rBV	89183	191392	19.99%	0.903%
16	8.183	1158	1164	1169	rVV	59385	110034	11.49%	0.519%
17	8.269	1173	1178	1182	rBV	131520	219580	22.93%	1.036%
18	8.305	1182	1184	1190	rVB	57378	87151	9.10%	0.411%
19	8.433	1200	1205	1214	rBV	143858	254110	26.54%	1.199%
20	8.610	1226	1234	1236	rBV4	87751	167009	17.44%	0.788%
21	8.647	1236	1240	1248	rVB	400785	653679	68.27%	3.085%
22	8.720	1248	1252	1256	rBV	26734	48022	5.02%	0.227%
23	8.909	1277	1283	1290	rBV2	99400	165481	17.28%	0.781%
24	9.098	1306	1314	1321	rBV4	29238	74365	7.77%	0.351%
25	9.195	1327	1330	1339	rVB3	24374	45753	4.78%	0.216%
26	9.287	1340	1345	1350	rBV	33429	48458	5.06%	0.229%
27	9.384	1355	1361	1370	rBV	52387	94632	9.88%	0.447%
28	9.494	1374	1379	1386	rVV2	100964	196933	20.57%	0.929%
29	9.646	1402	1404	1416	rVB5	24973	73019	7.63%	0.345%
30	9.756	1416	1422	1428	rBV3	36578	80355	8.39%	0.379%
31	9.817	1428	1432	1436	rVV2	40116	77799	8.13%	0.367%
32	9.866	1437	1440	1442	rVV2	39323	59355	6.20%	0.280%
33	9.903	1442	1446	1451	rVB	319346	436192	45.55%	2.058%
34	10.006	1458	1463	1467	rBV	243387	316563	33.06%	1.494%
35	10.055	1467	1471	1476	rVV	430526	585330	61.13%	2.762%
36	10.195	1489	1494	1499	rVB	88083	127818	13.35%	0.603%

LSC Area Percent Report

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Sample : 06017-01
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-7.5-8.5

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_X\Method\82X121323W.M
Title : SW846 8260

37	10.262	1499	1505	1509	rBV4	45135	117789	12.30%	0.556%
38	10.305	1509	1512	1516	rVB2	69523	103884	10.85%	0.490%
39	10.360	1516	1521	1531	rVB4	157032	319032	33.32%	1.505%
40	10.585	1553	1558	1562	rBV	84643	112468	11.75%	0.531%
41	10.671	1566	1572	1577	rBV3	70339	116126	12.13%	0.548%
42	10.756	1582	1586	1587	rBV2	35977	47317	4.94%	0.223%
43	10.780	1587	1590	1594	rVB	91926	109534	11.44%	0.517%
44	10.854	1594	1602	1606	rBV3	77242	143329	14.97%	0.676%
45	10.963	1613	1620	1627	rBV4	48469	119087	12.44%	0.562%
46	11.030	1627	1631	1634	rBV2	66321	95361	9.96%	0.450%
47	11.085	1634	1640	1642	rBV	460953	646165	67.48%	3.049%
48	11.110	1642	1644	1650	rVB2	256130	360328	37.63%	1.700%
49	11.189	1650	1657	1661	rBV	182705	237360	24.79%	1.120%
50	11.274	1665	1671	1673	rBV5	34397	76997	8.04%	0.363%
51	11.305	1673	1676	1683	rVB	155817	194881	20.35%	0.920%
52	11.433	1694	1697	1700	rBV	32523	43445	4.54%	0.205%
53	11.475	1700	1704	1711	rVB2	171498	248273	25.93%	1.172%
54	11.628	1723	1729	1734	rBV2	52482	97776	10.21%	0.461%
55	11.683	1734	1738	1740	rBV	48586	55006	5.74%	0.260%
56	11.713	1740	1743	1746	rVB	48960	45908	4.79%	0.217%
57	11.750	1746	1749	1759	rVB3	99880	205767	21.49%	0.971%
58	11.841	1759	1764	1766	rBV2	63202	93837	9.80%	0.443%
59	11.890	1770	1772	1775	rVB	48846	49128	5.13%	0.232%
60	11.933	1775	1779	1783	rBV2	62288	95415	9.96%	0.450%
61	12.024	1788	1794	1799	rBV2	466199	701907	73.31%	3.312%
62	12.073	1799	1802	1804	rBV2	68439	77141	8.06%	0.364%
63	12.097	1804	1806	1809	rVB	70332	71556	7.47%	0.338%
64	12.170	1816	1818	1825	rVB	109373	133860	13.98%	0.632%
65	12.244	1825	1830	1836	rBV	401670	570164	59.55%	2.691%
66	12.311	1836	1841	1849	rVB3	222878	464559	48.52%	2.192%
67	12.420	1854	1859	1863	rVB2	101717	140480	14.67%	0.663%
68	12.469	1863	1867	1872	rVB2	95978	148456	15.50%	0.701%
69	12.530	1873	1877	1883	rBV	100516	175580	18.34%	0.829%
70	12.615	1887	1891	1894	rBV	553920	643688	67.23%	3.037%
71	12.646	1894	1896	1900	rVB2	66316	67590	7.06%	0.319%
72	12.701	1900	1905	1914	rVB3	214463	419519	43.81%	1.980%
73	12.798	1916	1921	1926	rBV6	23097	47938	5.01%	0.226%
74	12.920	1933	1941	1945	rBV	421163	612171	63.93%	2.889%
75	12.963	1945	1948	1954	rVB	384746	496669	51.87%	2.344%
76	13.024	1954	1958	1960	rBV	118657	170063	17.76%	0.802%

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Integration Parameters: RTEINT.P
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Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
Title : SW846 8260

77	13.115	1968	1973	1975	rBV2	37965	52474	5.48%	0.248%
78	13.170	1978	1982	1986	rVV	322915	408526	42.67%	1.928%
79	13.213	1986	1989	1992	rVB	66174	71997	7.52%	0.340%
80	13.262	1992	1997	1998	rBV3	121441	158243	16.53%	0.747%
81	13.292	1998	2002	2008	rVV2	660495	957507	100.00%	4.518%
82	13.371	2012	2015	2020	rVB	131918	164425	17.17%	0.776%
83	13.420	2020	2023	2032	rVB4	96590	175856	18.37%	0.830%
84	13.506	2032	2037	2040	rBV2	93961	148386	15.50%	0.700%
85	13.542	2040	2043	2046	rVV	162748	207243	21.64%	0.978%
86	13.579	2046	2049	2056	rVV3	231679	396992	41.46%	1.873%
87	13.640	2056	2059	2060	rVV3	44162	52641	5.50%	0.248%
88	13.664	2060	2063	2069	rVV2	141037	252646	26.39%	1.192%
89	13.774	2077	2081	2087	rVB	539555	678354	70.85%	3.201%
90	13.926	2099	2106	2110	rBV3	120590	209888	21.92%	0.990%
91	13.969	2110	2113	2117	rVV	81673	107143	11.19%	0.506%
92	14.073	2126	2130	2135	rBV	173925	212554	22.20%	1.003%
93	14.213	2149	2153	2158	rBV2	169892	261096	27.27%	1.232%
94	14.322	2167	2171	2176	rBV	85149	122038	12.75%	0.576%
95	14.371	2176	2179	2183	rVB	112990	132111	13.80%	0.623%
96	14.481	2193	2197	2202	rBV4	27925	58461	6.11%	0.276%
97	14.633	2214	2222	2233	rVB	508406	747246	78.04%	3.526%
98	14.780	2241	2246	2250	rBV	248720	303313	31.68%	1.431%
99	15.475	2355	2360	2373	rVV	41718	74531	7.78%	0.352%
100	15.615	2377	2383	2386	rVV	43466	68256	7.13%	0.322%

Sum of corrected areas: 21191706

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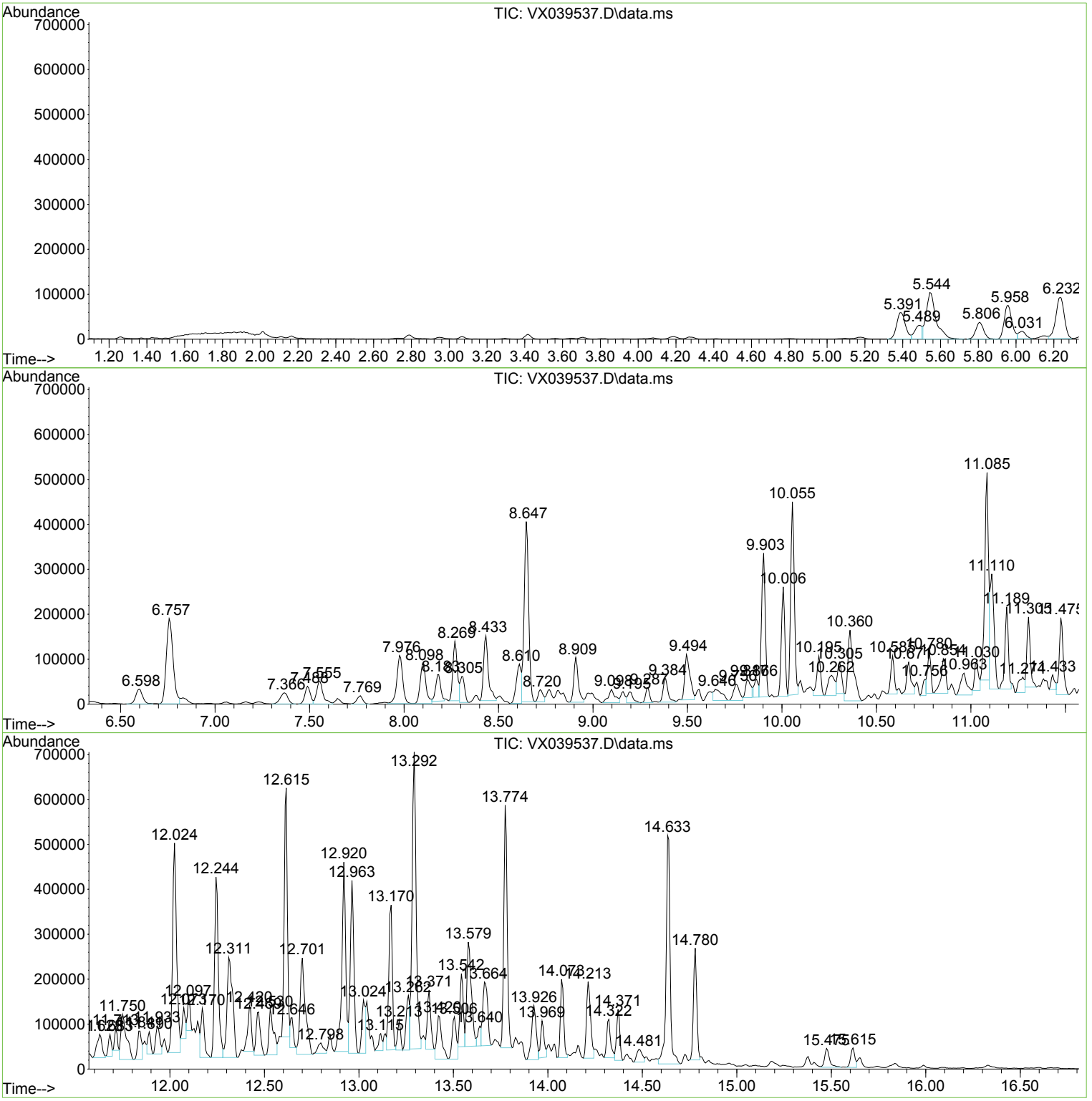
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-7.5-8.5

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122923\
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Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-7.5-8.5

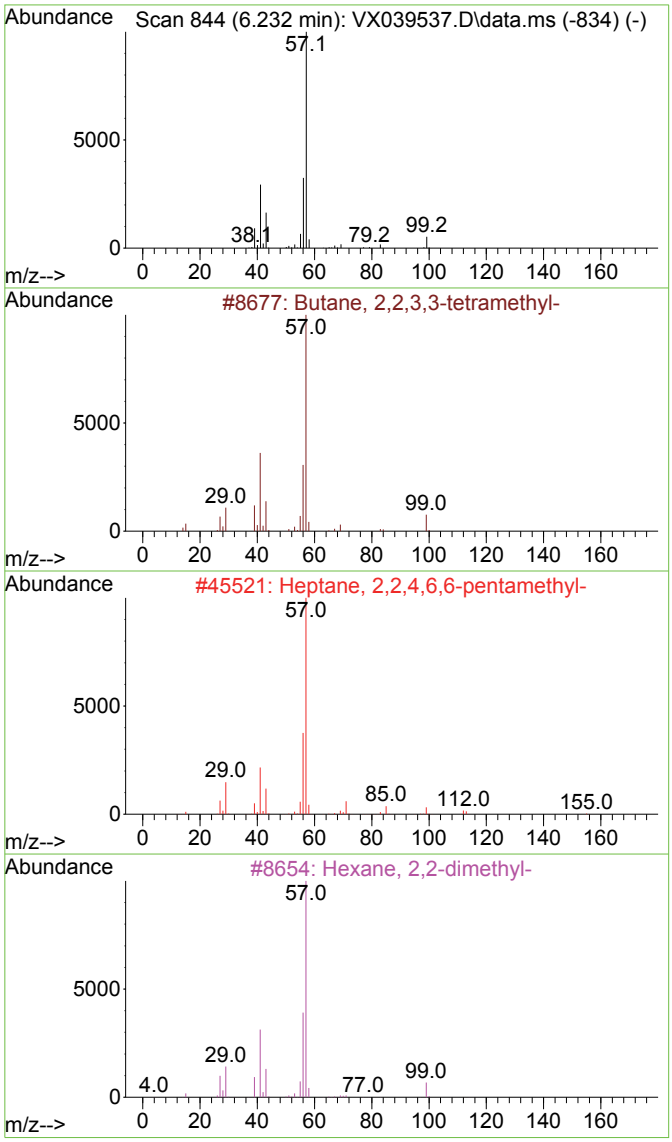
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
Quant Title : SW846 8260

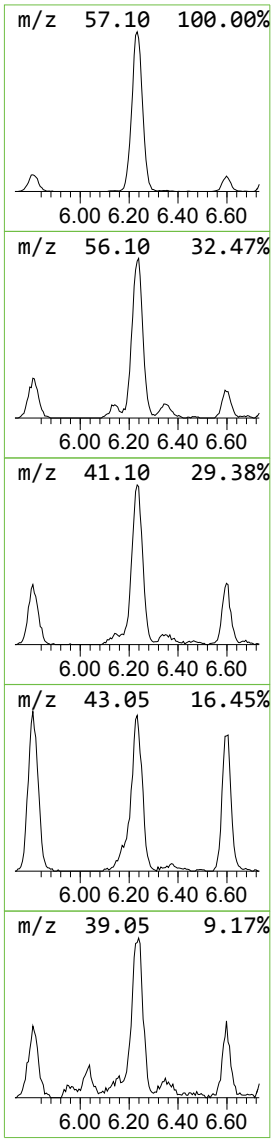
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Butane, 2,2,3,3-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.232	30.55 ug/l	287056	1,4-Difluorobenzene	6.757

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	72
2	Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	72
3	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	64
4	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	64
5	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122923\
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Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-7.5-8.5

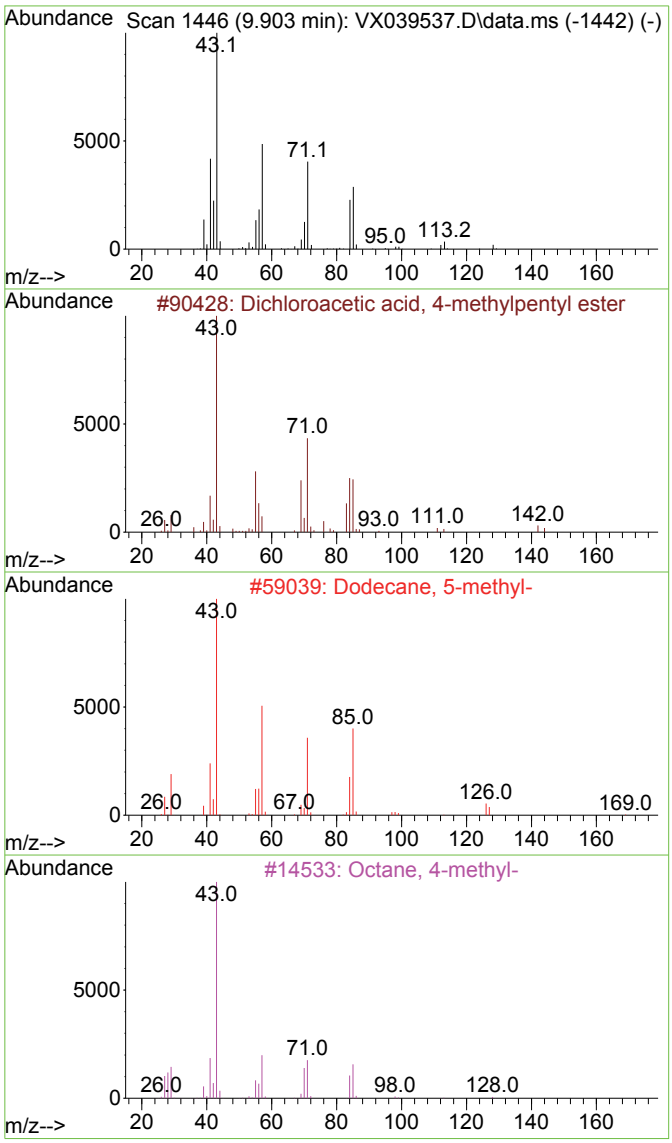
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
Quant Title : SW846 8260

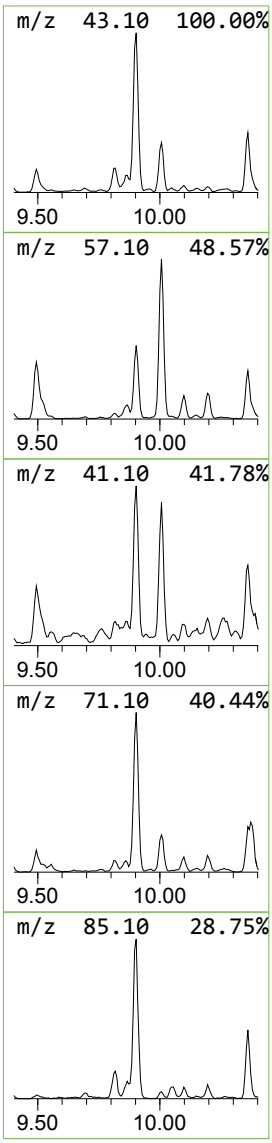
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Dichloroacetic acid, 4-meth... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.903	37.26 ug/l	436192	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dichloroacetic acid, 4-methylpen...	212	C8H14Cl2O2	1000282-43-7	64
2	Dodecane, 5-methyl-	184	C13H28	017453-93-9	59
3	Octane, 4-methyl-	128	C9H20	002216-34-4	58
4	Carbonic acid, nonyl prop-1-en-2...	228	C13H24O3	1000382-53-9	53
5	Heptane, 2,6-dimethyl-	128	C9H20	001072-05-5	53





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Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-7.5-8.5

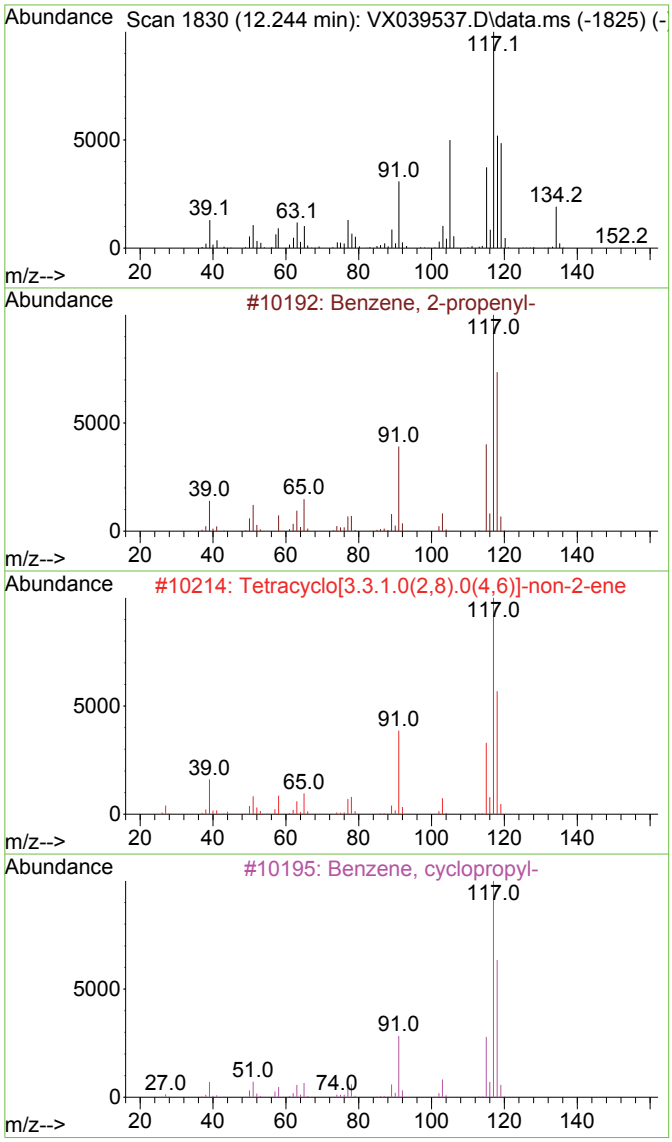
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
Quant Title : SW846 8260

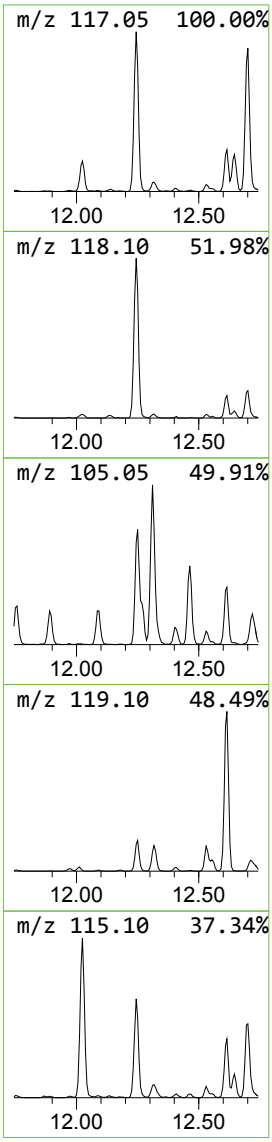
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 2-propenyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	40.62 ug/l	570164	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-propenyl-	118	C9H10	000300-57-2	91
2	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	60
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	60
4	Benzene, 1-propenyl-	118	C9H10	000637-50-3	60
5	Deltacyclene	118	C9H10	007785-10-6	55





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Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-7.5-8.5

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
Quant Title : SW846 8260

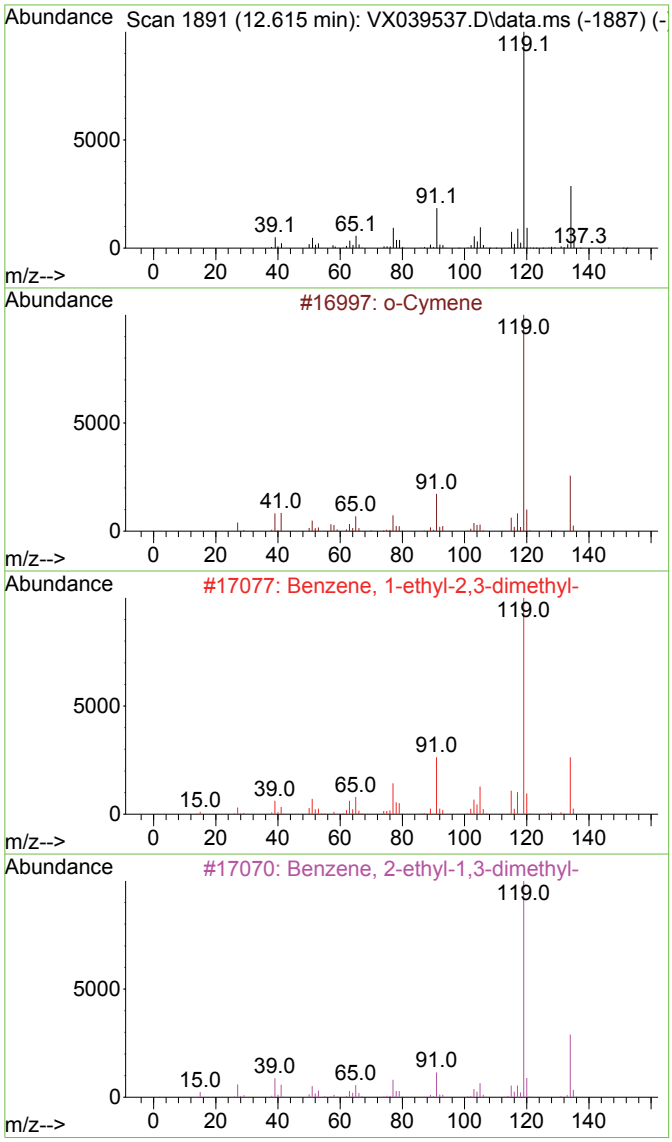
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

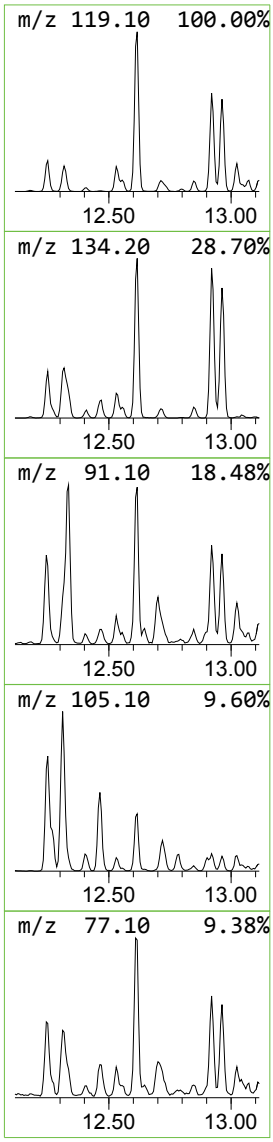
****

Peak Number 4 o-Cymene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.616	45.85 ug/l	643688	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	95
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
3	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	93





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122923\
Data File : VX039537.D
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Operator : JC/MD
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ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-7.5-8.5

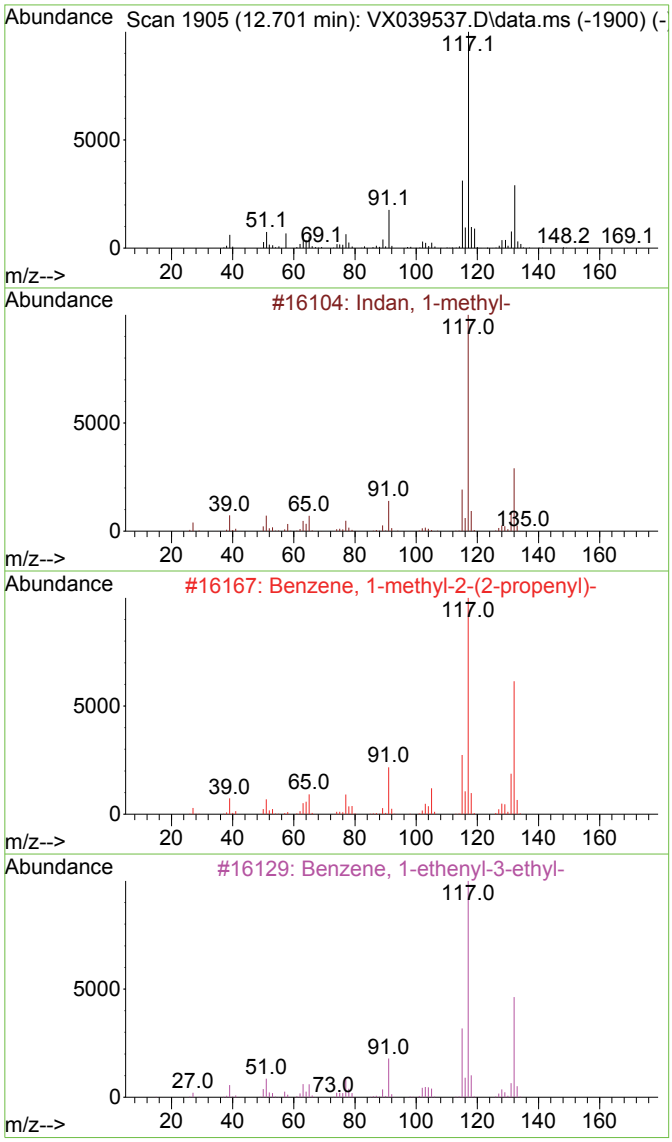
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
Quant Title : SW846 8260

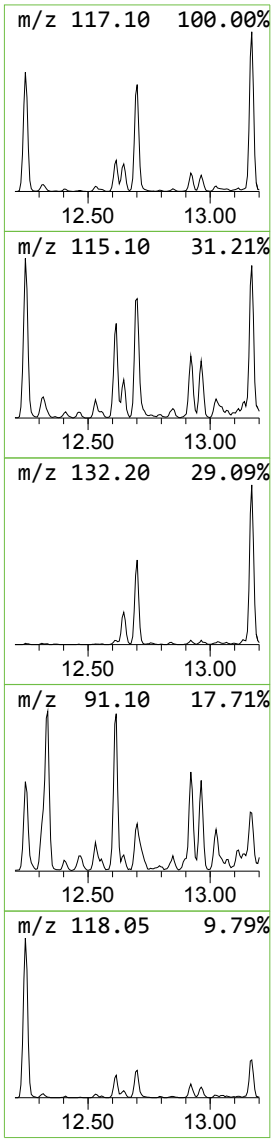
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Indan, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	29.88 ug/l	419519	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	87
2	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87
3	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
4	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	86





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Data File : VX039537.D
Acq On : 29 Dec 2023 12:04
Operator : JC/MD
Sample : 06017-01
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-7.5-8.5

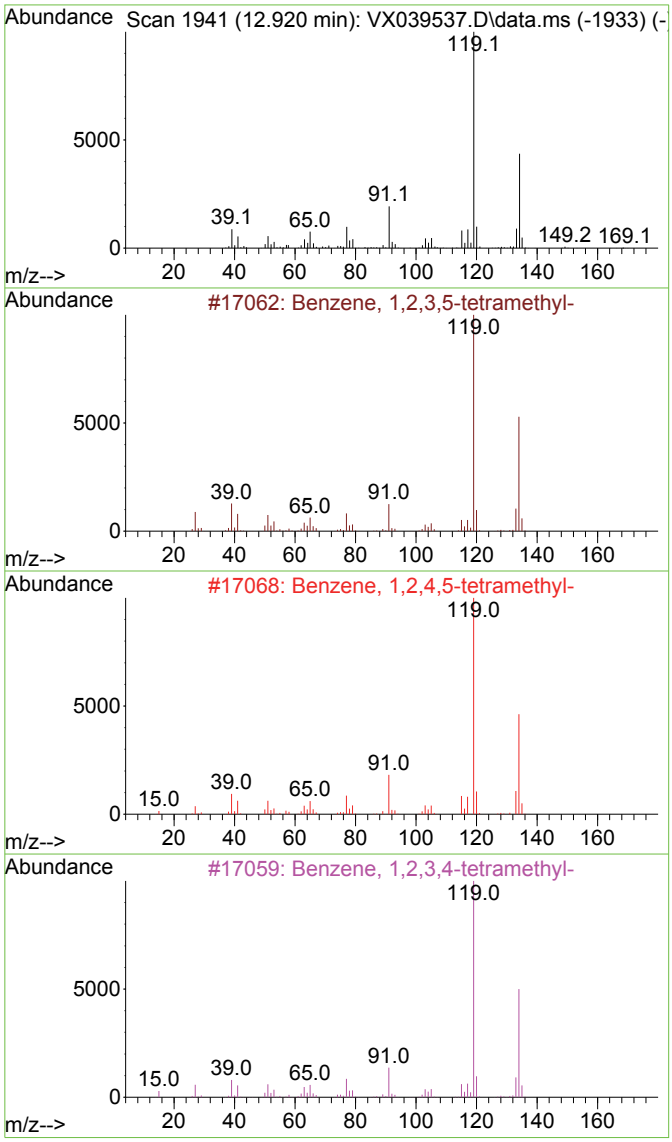
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
Quant Title : SW846 8260

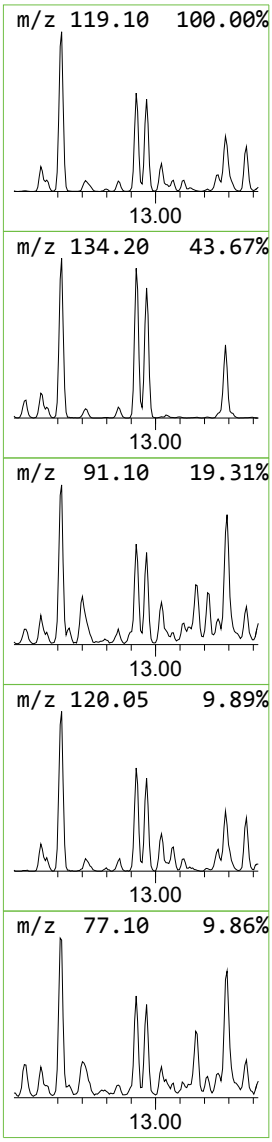
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.920	43.61 ug/l	612171	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4	o-Cymene	134	C10H14	000527-84-4	94
5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91





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Data File : VX039537.D
Acq On : 29 Dec 2023 12:04
Operator : JC/MD
Sample : 06017-01
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-7.5-8.5

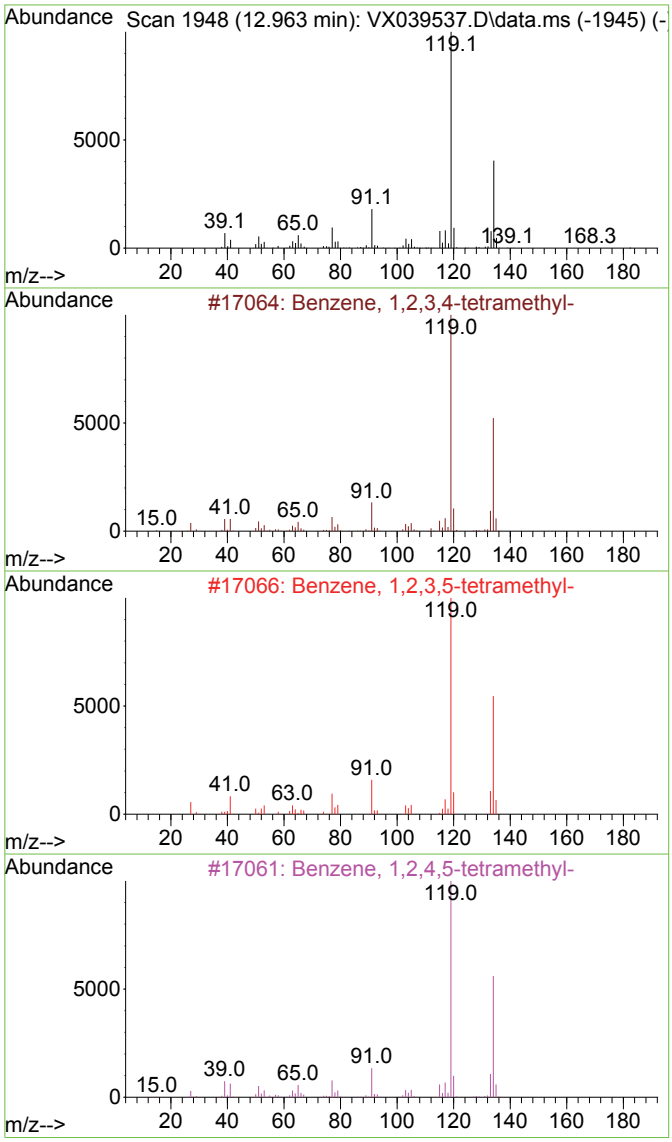
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
Quant Title : SW846 8260

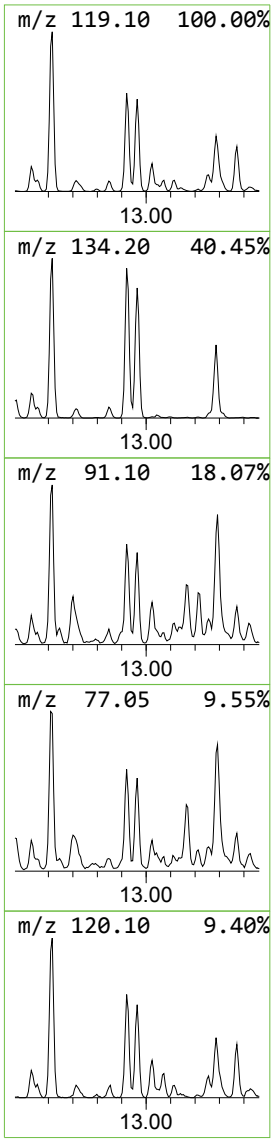
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.963	35.38 ug/l	496669	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122923\
Data File : VX039537.D
Acq On : 29 Dec 2023 12:04
Operator : JC/MD
Sample : 06017-01
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-7.5-8.5

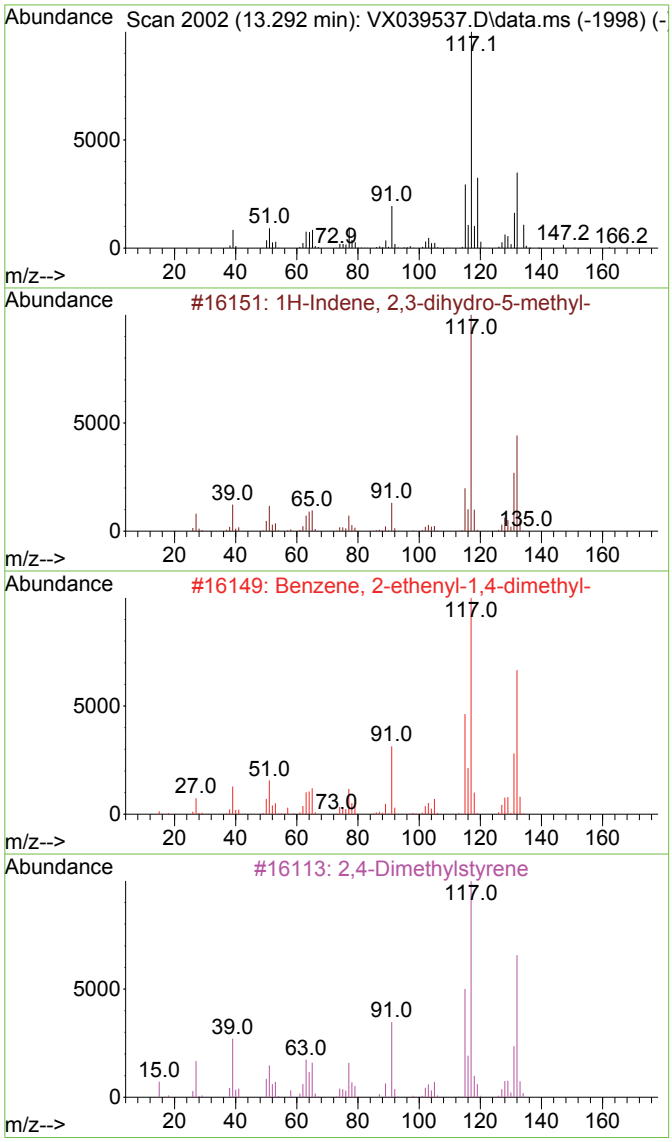
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
Quant Title : SW846 8260

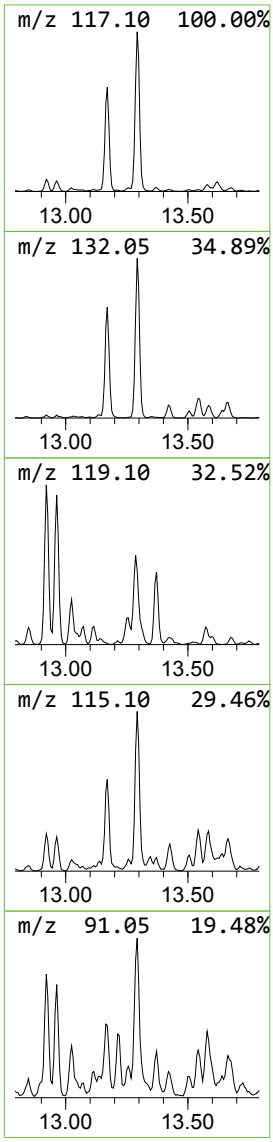
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	68.21 ug/l	957507	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
2	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
3	2,4-Dimethylstyrene	132	C10H12	002234-20-0	92
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	81
5	1-Phenyl-1-butene	132	C10H12	000824-90-8	81





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122923\
Data File : VX039537.D
Acq On : 29 Dec 2023 12:04
Operator : JC/MD
Sample : 06017-01
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-7.5-8.5

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
Quant Title : SW846 8260

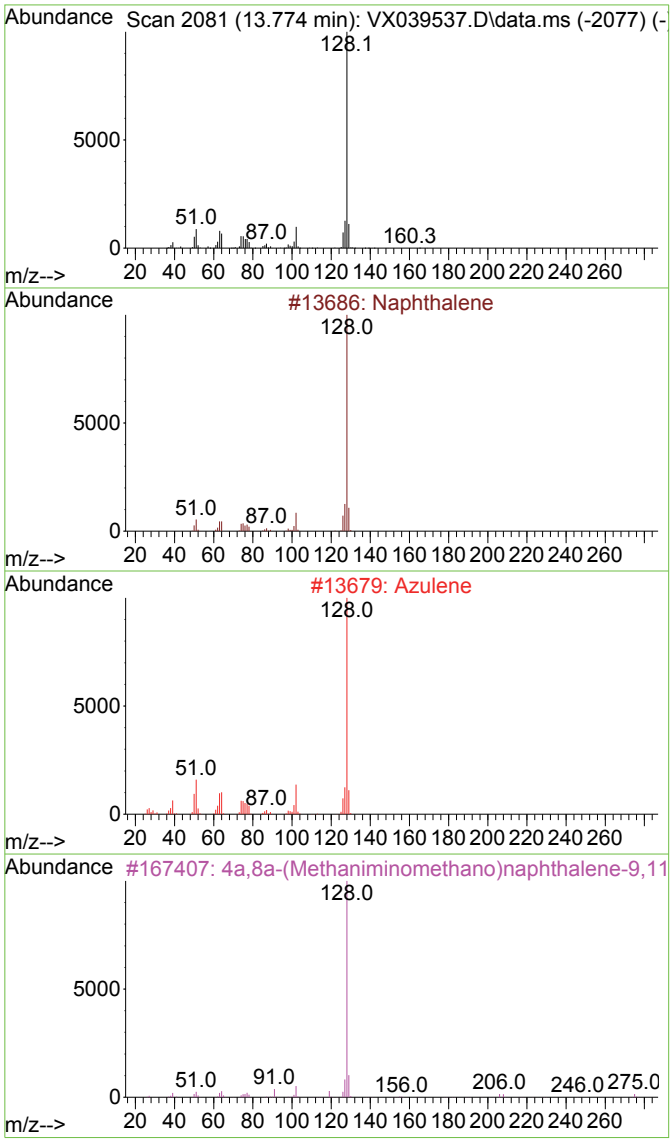
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

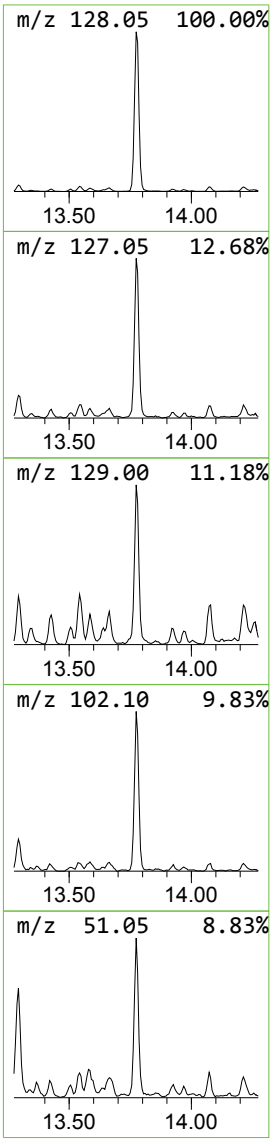
****

Peak Number 9 Azulene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.774	48.32 ug/l	678354	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene	128	C10H8	000091-20-3	95
2	Azulene	128	C10H8	000275-51-4	94
3	4a,8a-(Methaniminomethano)naphth...	275	C18H13NO2	069915-10-2	74
4	4-Phenylbut-3-ene-1-yne	128	C10H8	1000222-12-0	72
5	1,3-Dicyanobenzene	128	C8H4N2	000626-17-5	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122923\
Data File : VX039537.D
Acq On : 29 Dec 2023 12:04
Operator : JC/MD
Sample : 06017-01
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-7.5-8.5

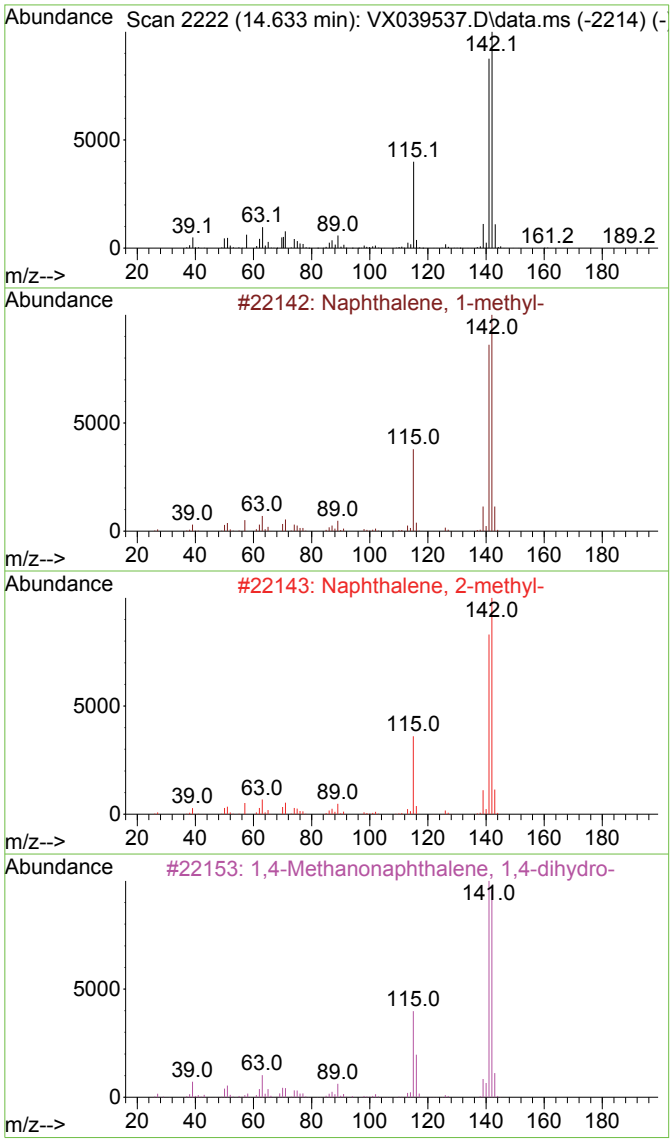
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
Quant Title : SW846 8260

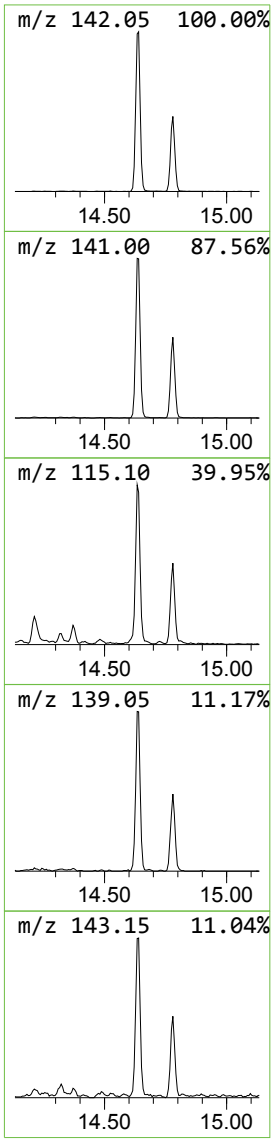
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 1,4-Methanonaphthalene, 1,4... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.633	53.23 ug/l	747246	1,4-Dichlorobenzene-d4	12.024

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	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	93
4	Benzocycloheptatriene	142	C11H10	000264-09-5	90
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	72





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122923\
Data File : VX039537.D
Acq On : 29 Dec 2023 12:04
Operator : JC/MD
Sample : 06017-01
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-7.5-8.5

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2,2,3,3...	6.232	30.6	ug/l	287056	2	6.757	469835	50.0
Dichloroacetic ...	9.903	37.3	ug/l	436192	3	10.055	585330	50.0
Benzene, 2-prop...	12.244	40.6	ug/l	570164	4	12.024	701907	50.0
o-Cymene	12.616	45.9	ug/l	643688	4	12.024	701907	50.0
Indan, 1-methyl-	12.701	29.9	ug/l	419519	4	12.024	701907	50.0
Benzene, 1,2,3,...	12.920	43.6	ug/l	612171	4	12.024	701907	50.0
Benzene, 1,2,3,...	12.963	35.4	ug/l	496669	4	12.024	701907	50.0
1H-Indene, 2,3-...	13.292	68.2	ug/l	957507	4	12.024	701907	50.0
Azulene	13.774	48.3	ug/l	678354	4	12.024	701907	50.0
1,4-Methanonaph...	14.633	53.2	ug/l	747246	4	12.024	701907	50.0