LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042823\
Data File : VX035394.D
Acq On : 28 Apr 2023 16:29
Operator : JC/MD
Sample : 02386-14 10X
Misc : 3.62g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SPA-4WC-23-04

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\82X042523W.M
Title : SW846 8260

Signal : TIC: VX035394.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.385	692	705	717	rBV	143072	415149	17.75%	1.131%
2	5.550	719	732	752	rVB	262891	766159	32.76%	2.087%
3	5.952	789	798	809	rBV	155422	416089	17.79%	1.133%
4	6.757	921	930	947	rBV	397715	963288	41.19%	2.624%
5	7.379	1021	1032	1043	rVB3	24959	57082	2.44%	0.155%
6	8.275	1173	1179	1183	rBV	27994	48105	2.06%	0.131%
7	8.433	1200	1205	1215	rVB	25568	47984	2.05%	0.131%
8	8.598	1226	1232	1235	rBV2	33584	60761	2.60%	0.165%
9	8.647	1235	1240	1247	rVV	835209	1322845	56.56%	3.603%
10	8.720	1247	1252	1257	rVB	96795	147245	6.30%	0.401%
11	8.915	1278	1284	1289	rBV	57797	89360	3.82%	0.243%
12	9.385	1351	1361	1371	rVB	24760	41815	1.79%	0.114%
13	9.500	1374	1380	1386	rBV3	35235	72545	3.10%	0.198%
14	9.561	1386	1390	1395	rVB2	38712	55711	2.38%	0.152%
15	9.823	1426	1433	1438	rBV3	34083	70148	3.00%	0.191%
16	9.903	1442	1446	1452	rVB	92596	128644	5.50%	0.350%
17	10.006	1457	1463	1466	rBV	73581	102434	4.38%	0.279%
18	10.055	1466	1471	1486	rVB2	833844	1377246	58.89%	3.751%
19	10.195	1487	1494	1499	rBV	795153	1062999	45.45%	2.895%
20	10.299	1503	1511	1517	rBV	1739559	2338658	100.00%	6.370%
21	10.360	1517	1521	1532	rVB2	162257	243298	10.40%	0.663%
22	10.586	1551	1558	1562	rBV2	99124	154409	6.60%	0.421%
23	10.640	1562	1567	1571	rBV	471480	604783	25.86%	1.647%
24	10.781	1586	1590	1594	rVB	195913	244124	10.44%	0.665%
25	10.854	1594	1602	1606	rBV2	156147	282537	12.08%	0.770%
26	10.896	1606	1609	1613	rVB2	78961	105955	4.53%	0.289%
27	10.957	1613	1619	1623	rBV2	120190	245430	10.49%	0.668%
28	11.031	1627	1631	1634	rVV2	103328	138913	5.94%	0.378%
29	11.079	1634	1639	1643	rVV2	879372	1252957	53.58%	3.413%
30	11.110	1643	1644	1650	rVB2	223366	249069	10.65%	0.678%
31	11.189	1650	1657	1660	rBV	204664	324581	13.88%	0.884%
32	11.220	1660	1662	1666	rVB3	95176	100117	4.28%	0.273%
33	11.305	1672	1676	1679	rBV	178748	208015	8.89%	0.567%
34	11.378	1684	1688	1694	rVV	624233	1051475	44.96%	2.864%
35	11.451	1694	1700	1702	rVV2	410393	680804	29.11%	1.854%
36	11.476	1702	1704	1711	rVB	548079	642942	27.49%	1.751%

LSC Area Percent Report

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Instrument :
MSVOA_X
ClientSampleId :
SPA-4WC-23-04

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\82X042523W.M
Title : SW846 8260

37	11.610	1722	1726	1734	rVV2	254106	453928	19.41%	1.236%
38	11.683	1734	1738	1740	rVV2	106435	132599	5.67%	0.361%
39	11.713	1740	1743	1746	rVV	364345	451912	19.32%	1.231%
40	11.750	1746	1749	1759	rVB	1264531	1700284	72.70%	4.631%
41	11.835	1759	1763	1766	rBV	104441	166325	7.11%	0.453%
42	11.866	1766	1768	1770	rVV	113027	143274	6.13%	0.390%
43	11.890	1770	1772	1776	rVV3	157889	199179	8.52%	0.543%
44	11.939	1776	1780	1783	rVV2	287453	479077	20.49%	1.305%
45	11.969	1783	1785	1789	rVV3	439031	523907	22.40%	1.427%
46	12.018	1789	1793	1800	rVV3	983865	2066871	88.38%	5.630%
47	12.085	1800	1804	1809	rVV3	387288	779385	33.33%	2.123%
48	12.152	1812	1815	1816	rVV	108873	137241	5.87%	0.374%
49	12.171	1816	1818	1826	rVB2	161286	227579	9.73%	0.620%
50	12.244	1826	1830	1832	rBV2	192081	281924	12.05%	0.768%
51	12.268	1832	1834	1837	rVV	246526	257923	11.03%	0.703%
52	12.329	1837	1844	1849	rVV3	660294	1425463	60.95%	3.883%
53	12.421	1850	1859	1863	rVV	391878	544333	23.28%	1.483%
54	12.463	1863	1866	1872	rVB2	193333	295914	12.65%	0.806%
55	12.530	1872	1877	1879	rBV	225942	308018	13.17%	0.839%
56	12.555	1879	1881	1885	rVV	261190	325703	13.93%	0.887%
57	12.610	1887	1890	1894	rVV	319248	415912	17.78%	1.133%
58	12.640	1894	1895	1900	rVV2	56290	55800	2.39%	0.152%
59	12.701	1900	1905	1913	rVV2	194619	432208	18.48%	1.177%
60	12.799	1916	1921	1922	rBV3	50381	68594	2.93%	0.187%
61	12.847	1926	1929	1932	rVB2	100094	115291	4.93%	0.314%
62	12.890	1932	1936	1938	rBV3	91331	129460	5.54%	0.353%
63	12.920	1938	1941	1944	rVV	236795	282488	12.08%	0.769%
64	12.963	1944	1948	1954	rVB	367429	496669	21.24%	1.353%
65	13.024	1954	1958	1959	rBV	68538	76314	3.26%	0.208%
66	13.042	1959	1961	1964	rVV	62570	82958	3.55%	0.226%
67	13.073	1964	1966	1969	rVB	65429	56598	2.42%	0.154%
68	13.164	1975	1981	1985	rVB3	116689	183973	7.87%	0.501%
69	13.213	1985	1989	1992	rVB	55103	61187	2.62%	0.167%
70	13.262	1992	1997	1998	rBV2	98097	158040	6.76%	0.430%
71	13.292	1998	2002	2007	rVB2	279347	436911	18.68%	1.190%
72	13.372	2012	2015	2019	rVV2	81409	108223	4.63%	0.295%
73	13.420	2019	2023	2031	rVB	248302	350824	15.00%	0.956%
74	13.585	2045	2050	2057	rBV3	289318	454891	19.45%	1.239%
75	13.676	2060	2065	2068	rVB3	67625	87537	3.74%	0.238%
76	13.774	2077	2081	2088	rVB	363168	484962	20.74%	1.321%

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ClientSampleId :
SPA-4WC-23-04

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

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Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Title : SW846 8260

77	13.847	2088	2093	2096	rBV6	28688	51270	2.19%	0.140%
78	13.932	2102	2107	2111	rVB3	35203	58366	2.50%	0.159%
79	14.073	2127	2130	2134	rBV2	42418	58176	2.49%	0.158%
80	14.213	2149	2153	2161	rVB2	57276	99845	4.27%	0.272%
81	14.323	2166	2171	2175	rVV	42888	63899	2.73%	0.174%
82	14.487	2193	2198	2206	rBV7	28113	81042	3.47%	0.221%
83	14.634	2215	2222	2226	rBV	394607	559448	23.92%	1.524%
84	14.780	2239	2246	2254	rVV	225082	334342	14.30%	0.911%
85	15.133	2294	2304	2313	rVV	469510	1337117	57.17%	3.642%
86	15.207	2313	2316	2321	rVB	163054	220569	9.43%	0.601%
87	15.322	2330	2335	2340	rBV	374870	515297	22.03%	1.404%
88	15.377	2340	2344	2347	rVV2	75426	107231	4.59%	0.292%
89	15.414	2347	2350	2355	rVB5	55635	72442	3.10%	0.197%
90	15.475	2355	2360	2365	rBV	182058	299246	12.80%	0.815%
91	15.609	2377	2382	2386	rVV2	197434	357838	15.30%	0.975%
92	15.652	2386	2389	2397	rVB	134899	189504	8.10%	0.516%
93	15.835	2411	2419	2420	rBV2	53078	108812	4.65%	0.296%
94	15.987	2440	2444	2454	rVB5	38649	81417	3.48%	0.222%
95	16.091	2454	2461	2472	rVV3	80112	199639	8.54%	0.544%
96	16.188	2474	2477	2483	rVB3	32505	63849	2.73%	0.174%
97	16.274	2486	2491	2499	rBV9	21902	67931	2.90%	0.185%
98	16.353	2499	2504	2512	rVB2	38349	71619	3.06%	0.195%
99	16.597	2540	2544	2549	rVB2	44996	77004	3.29%	0.210%
100	16.658	2549	2554	2565	rVB2	39108	85264	3.65%	0.232%

Sum of corrected areas: 36714522

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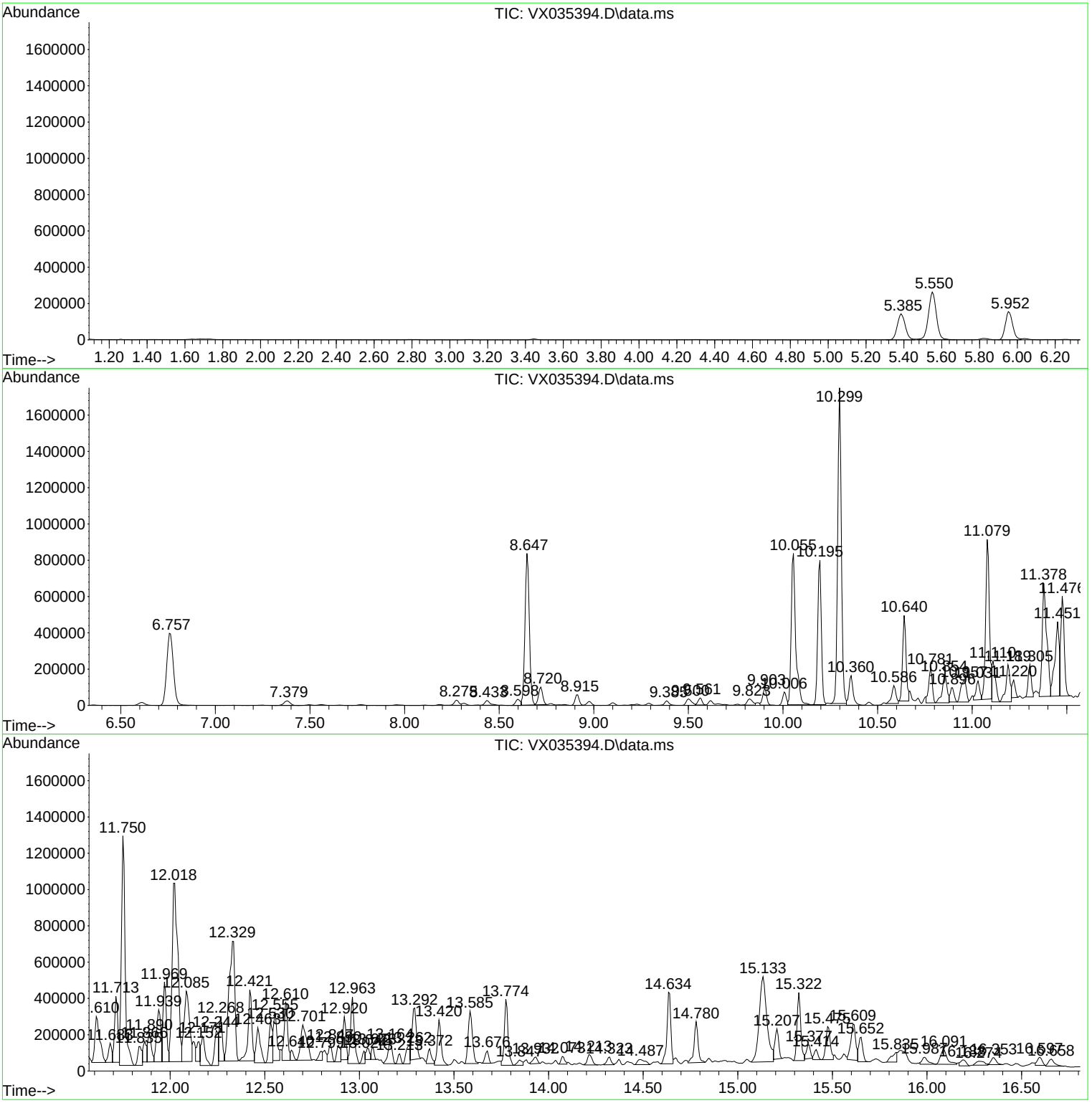
Instrument :
MSVOA_X
ClientSampleId :
SPA-4WC-23-04

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042823\
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Instrument :
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ClientSampleId :
SPA-4WC-23-04

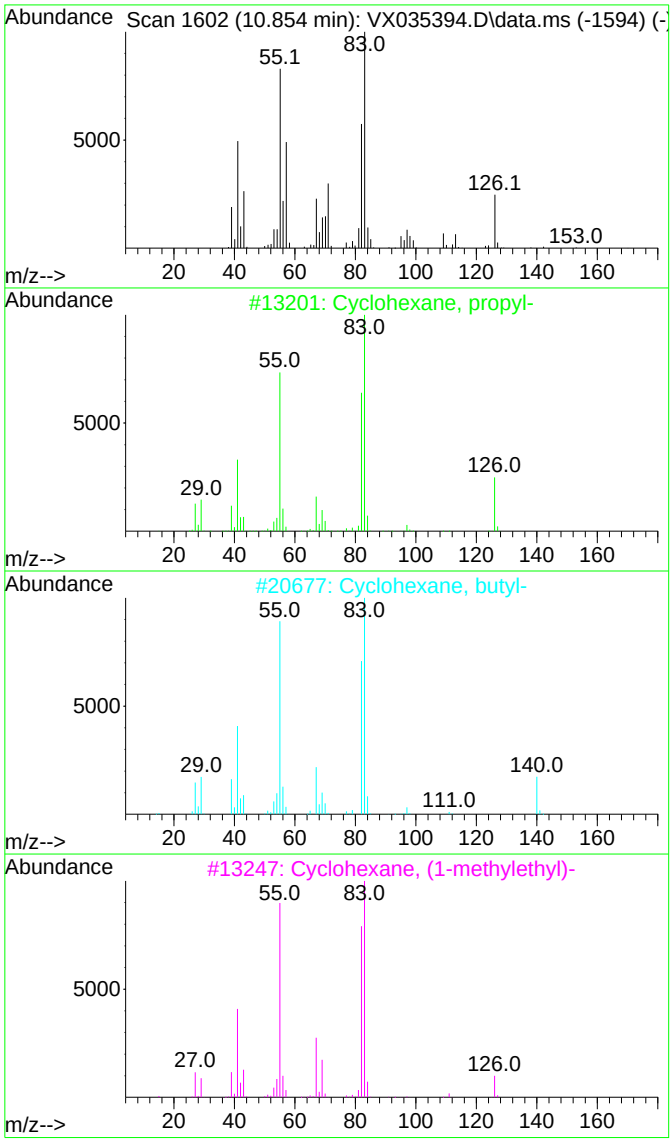
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

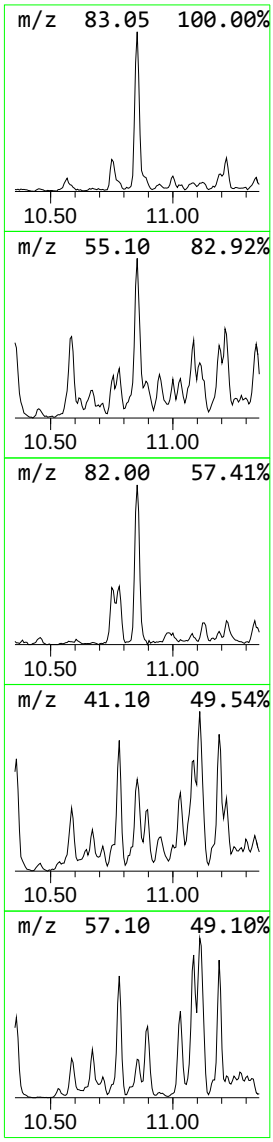
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Cyclohexane, propyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.854	10.26 ug/l	282537	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, propyl-	126	C9H18	001678-92-8	81
2	Cyclohexane, butyl-	140	C10H20	001678-93-9	59
3	Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	58
4	Cyclohexane, ethyl-	112	C8H16	001678-91-7	53
5	Cyclohexane, 1,1'-(1,4-butanediy...	222	C16H30	006165-44-2	53





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Instrument :
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ClientSampleId :
SPA-4WC-23-04

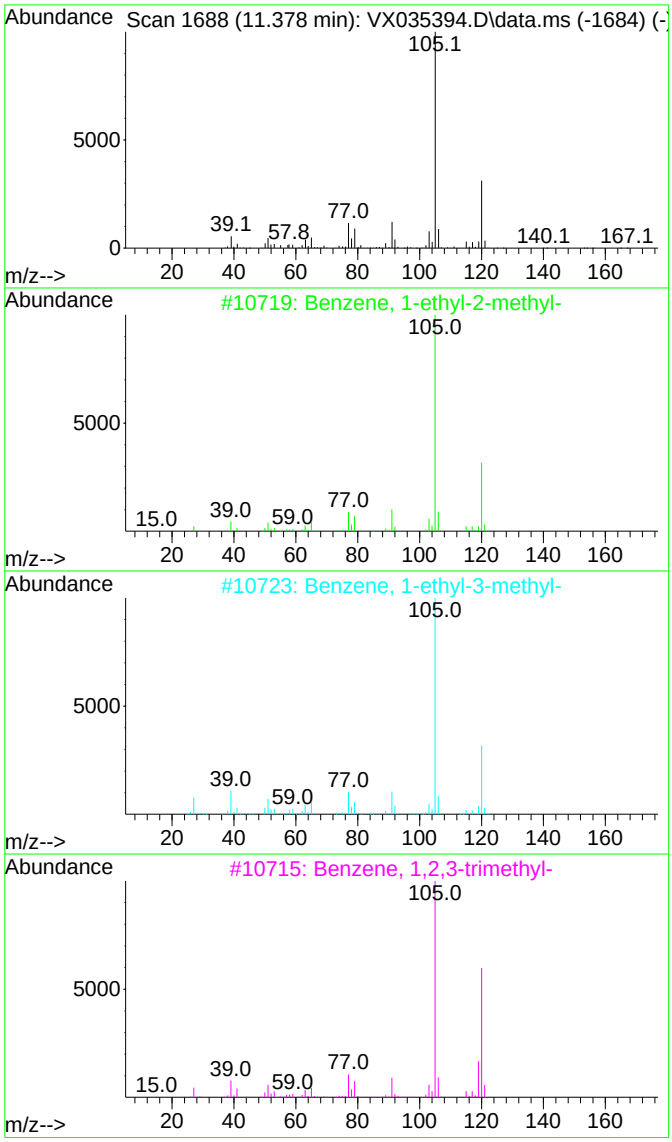
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

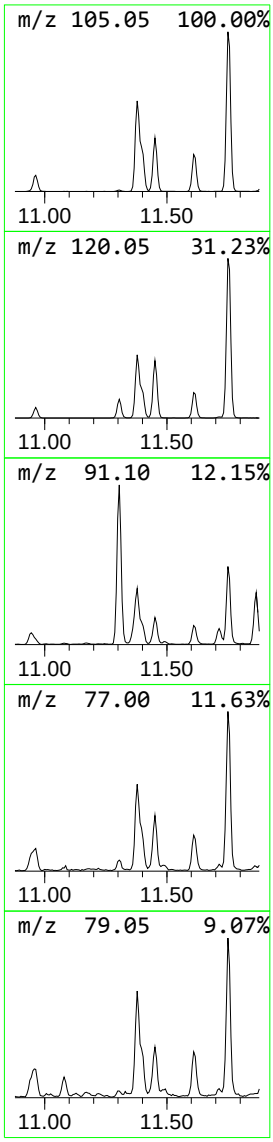
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.378	25.44 ug/l	1051480	1,4-Dichlorobenzene-d4	12.024

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, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





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Instrument :
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ClientSampleId :
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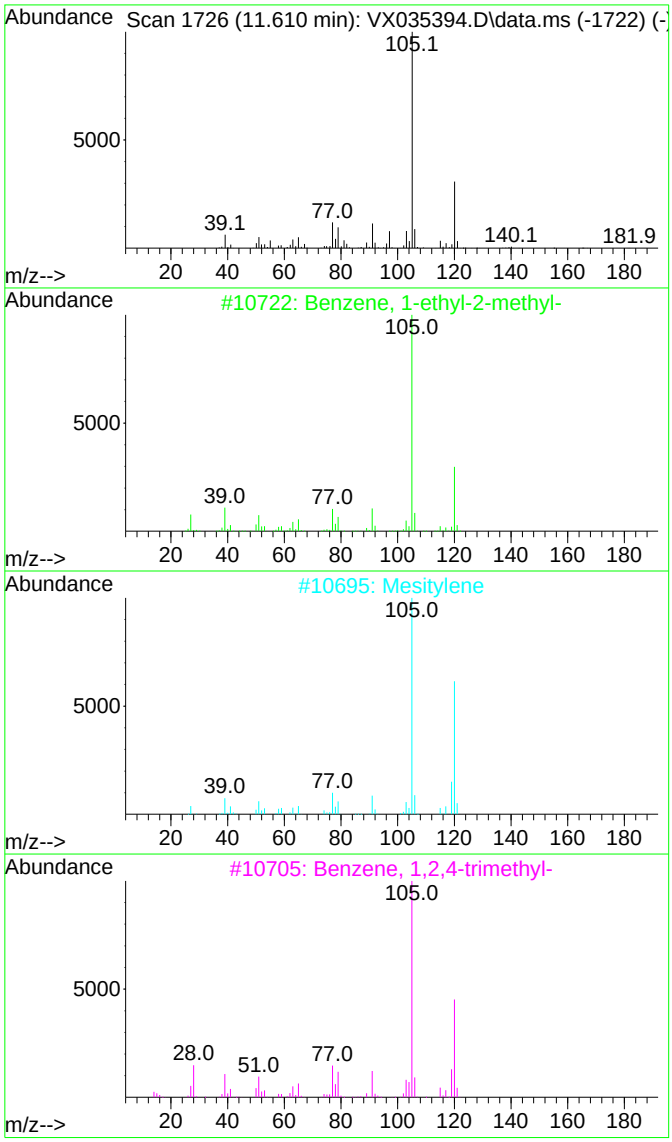
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

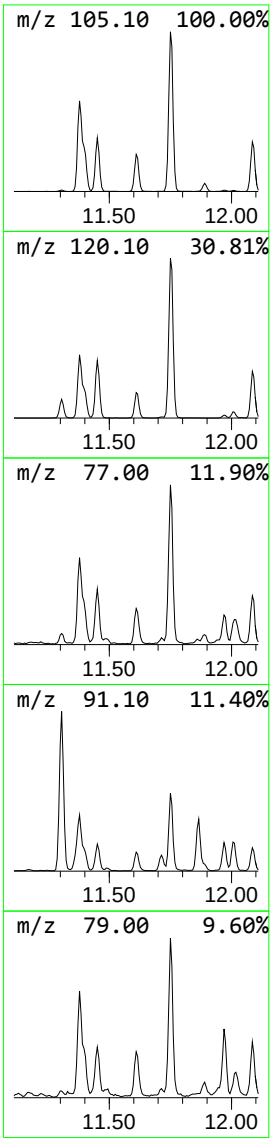
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1-ethyl-3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.610	10.98 ug/l	453928	1,4-Dichlorobenzene-d4	12.024

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	90
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90





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Instrument :
MSVOA_X
ClientSampleId :
SPA-4WC-23-04

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

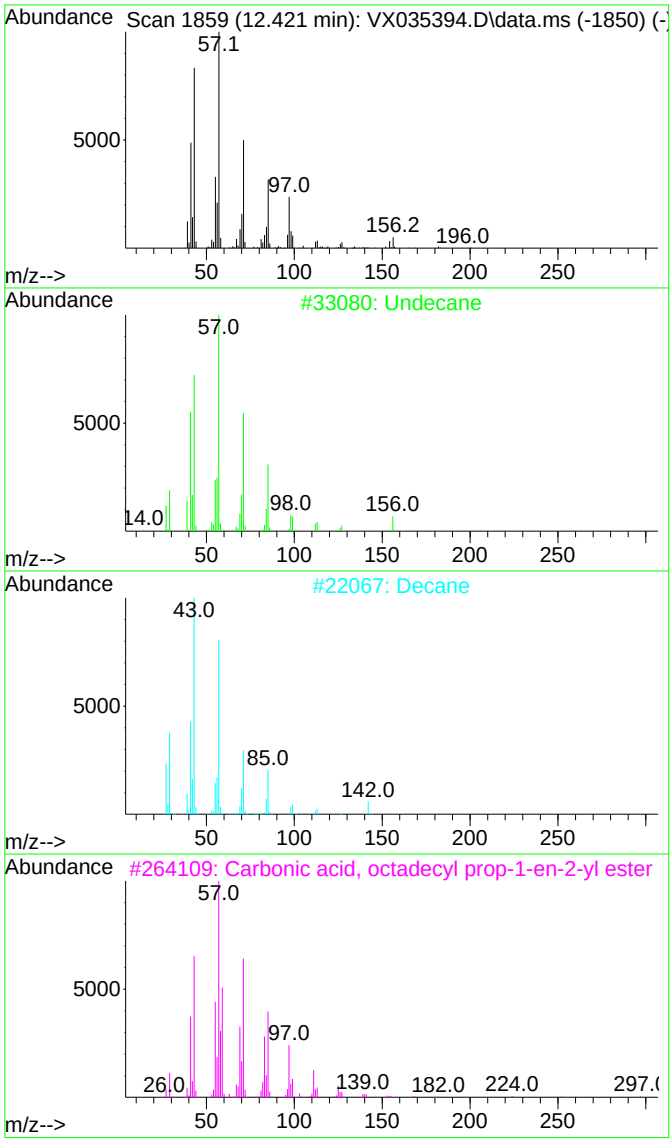
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

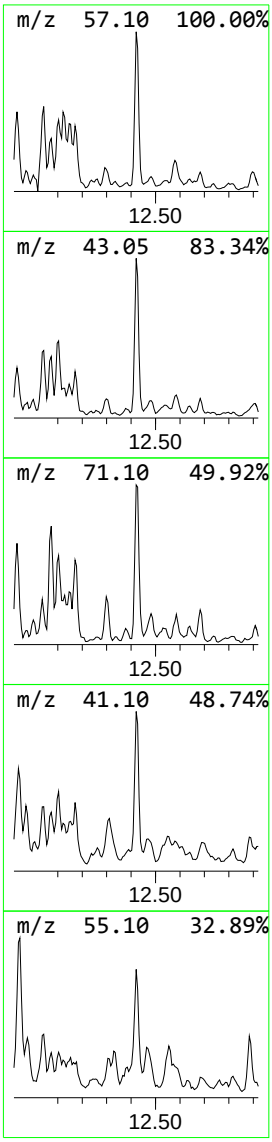
****

Peak Number 4 Undecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.421	13.17 ug/l	544333	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane	156	C11H24	001120-21-4	96
2	Decane	142	C10H22	000124-18-5	90
3	Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	86
4	Carbonic acid, prop-1-en-2-yl te...	298	C18H34O3	1000382-90-9	80
5	Tetradecane	198	C14H30	000629-59-4	72





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MSVOA_X
ClientSampleId :
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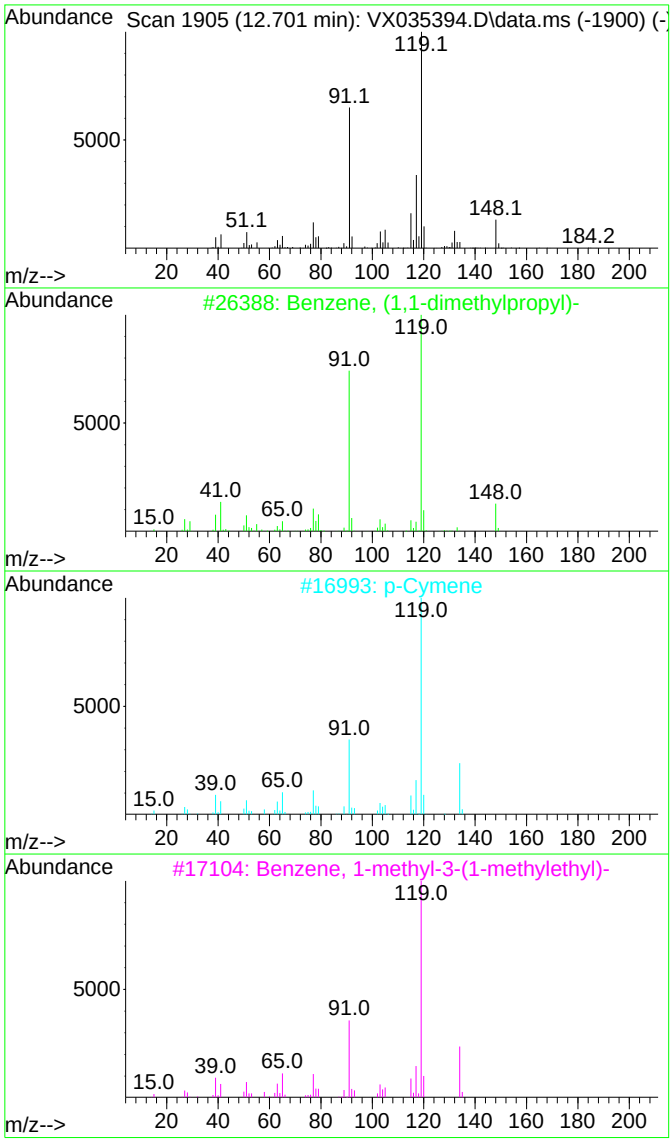
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

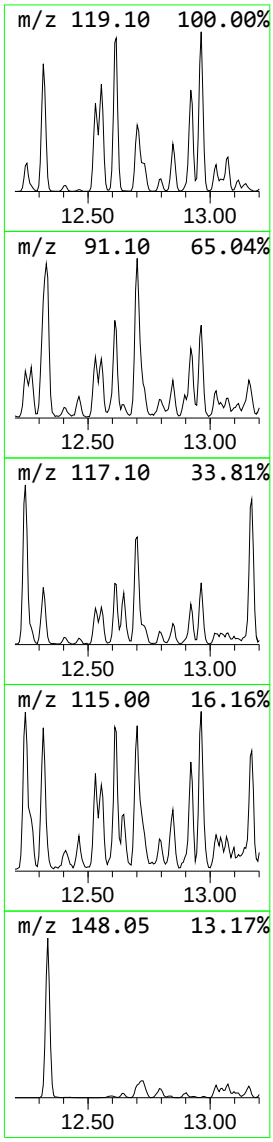
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, (1,1-dimethylpropyl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	10.46 ug/l	432208	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	81
2	p-Cymene	134	C10H14	000099-87-6	74
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	72
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	59
5	(6-Chloro-2-methylhexan-2-yl)ben...	210	C13H19Cl	1417621-45-4	59





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Data File : VX035394.D
Acq On : 28 Apr 2023 16:29
Operator : JC/MD
Sample : 02386-14 10X
Misc : 3.62g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SPA-4WC-23-04

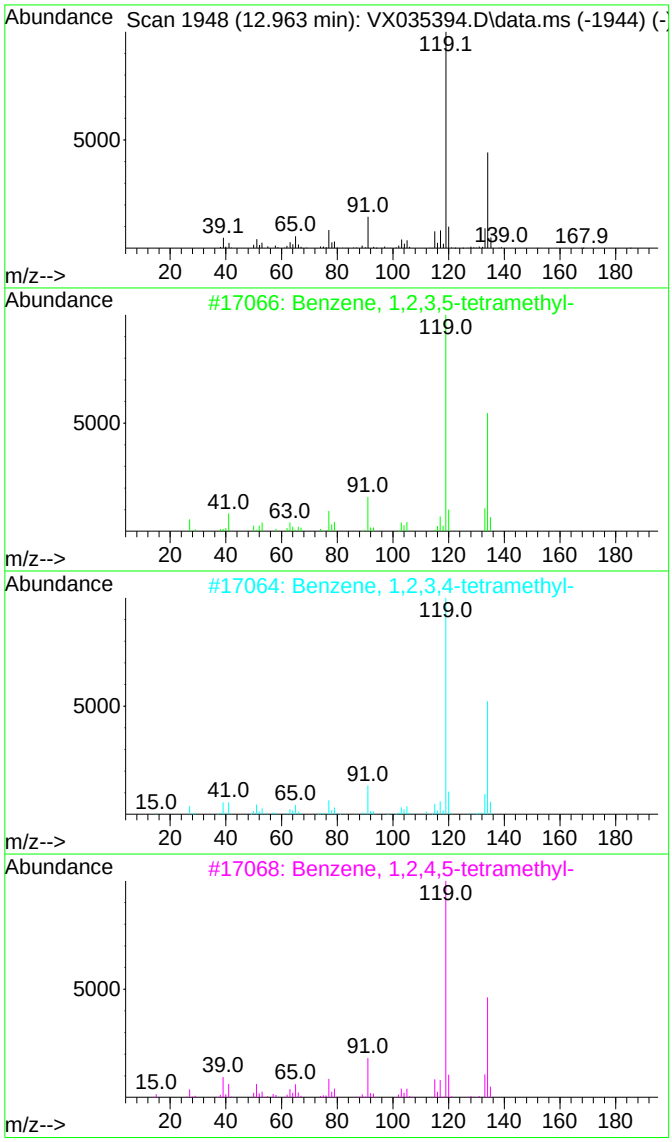
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

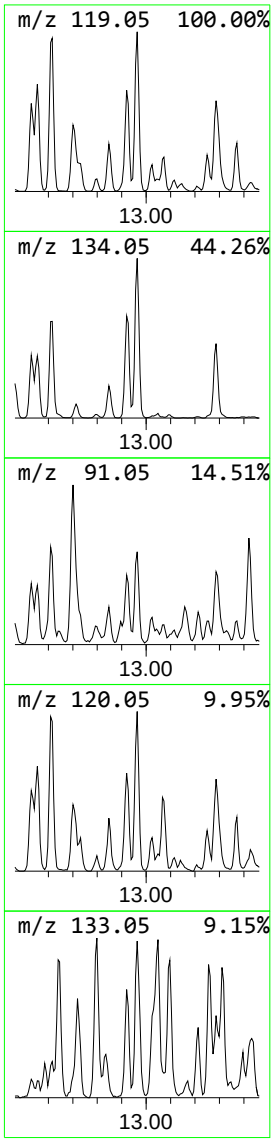
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.963	12.02 ug/l	496669	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	97
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4	o-Cymene	134	C10H14	000527-84-4	94
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93





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Data File : VX035394.D
Acq On : 28 Apr 2023 16:29
Operator : JC/MD
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ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SPA-4WC-23-04

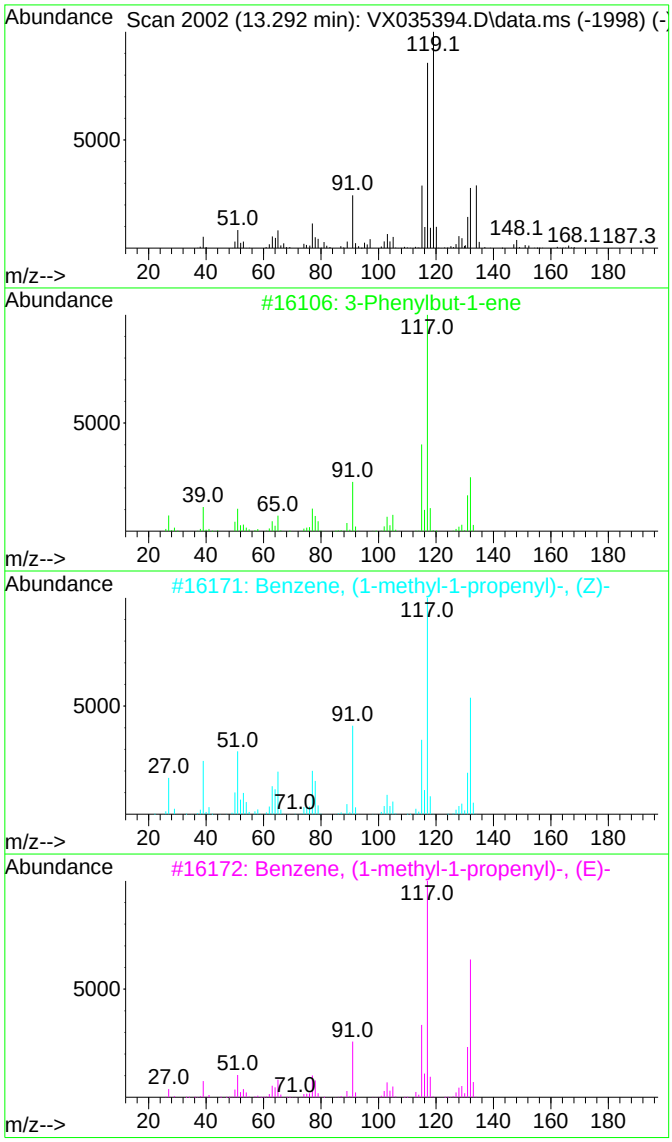
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

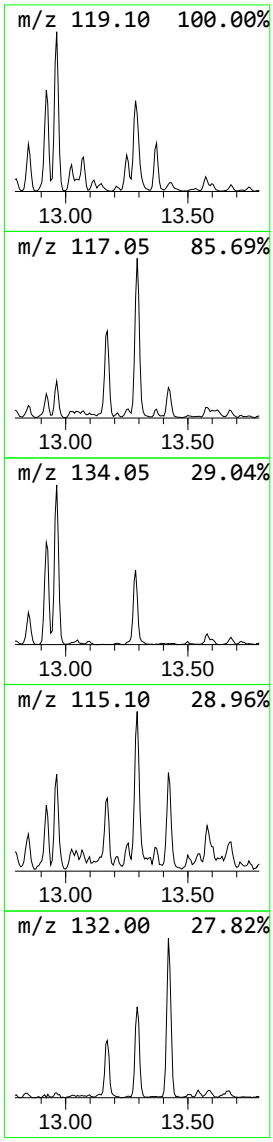
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 3-Phenylbut-1-ene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	10.57 ug/l	436911	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
2	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	70
3	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	55
4	Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	55
5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	55





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042823\
Data File : VX035394.D
Acq On : 28 Apr 2023 16:29
Operator : JC/MD
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Misc : 3.62g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SPA-4WC-23-04

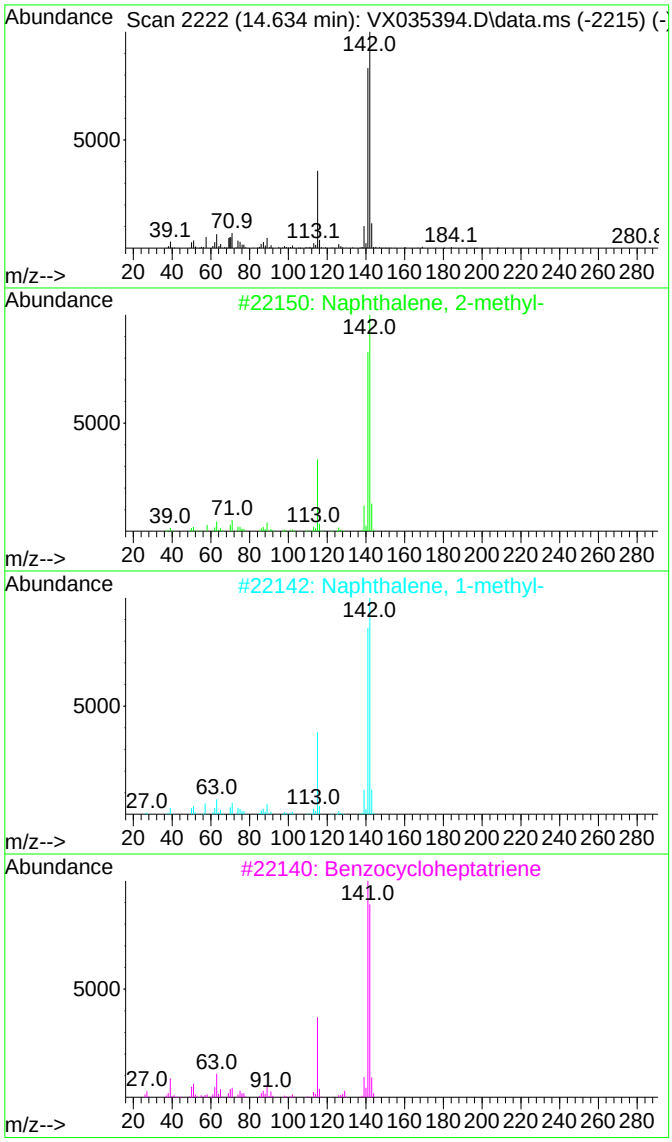
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

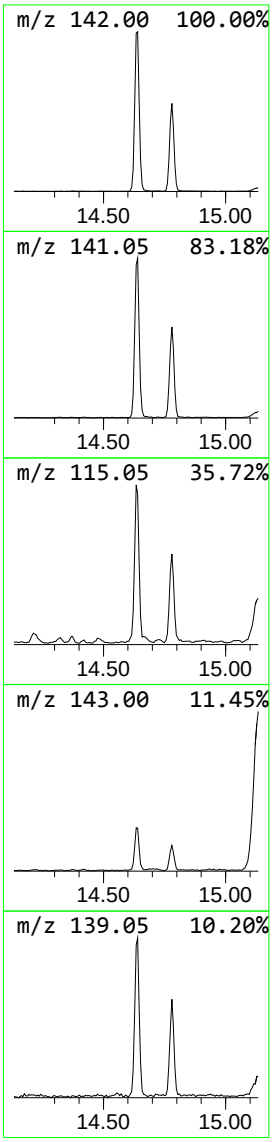
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzocycloheptatriene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.634	13.53 ug/l	559448	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3	Benzocycloheptatriene	142	C11H10	000264-09-5	91
4	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	90
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64





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Operator : JC/MD
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Misc : 3.62g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SPA-4WC-23-04

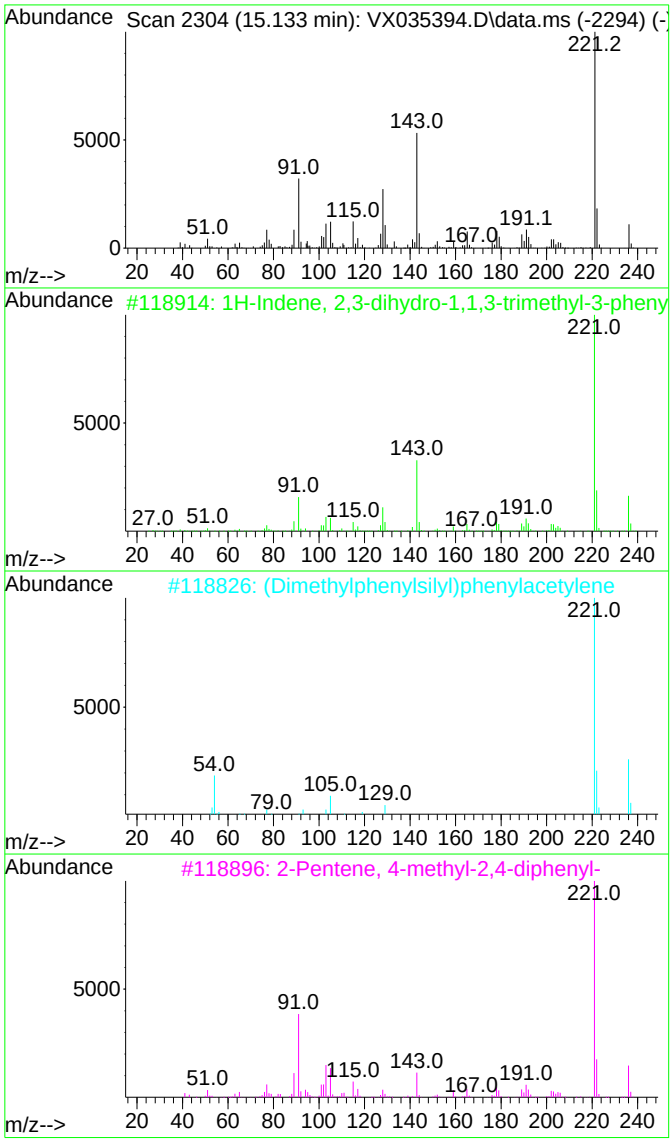
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

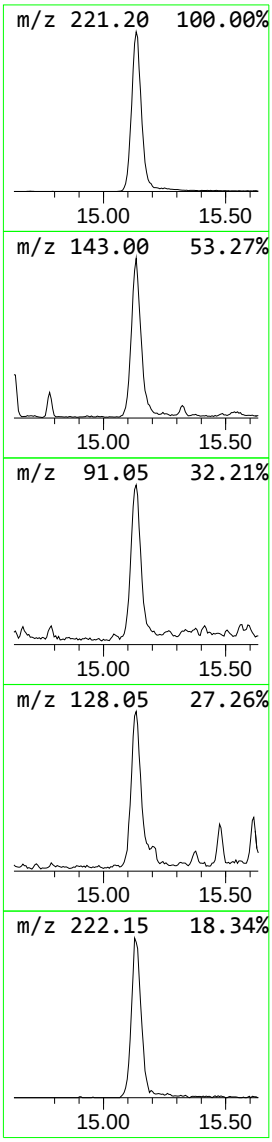
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 1H-Indene, 2,3-dihydro-1,1,... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.134	32.35 ug/l	1337120	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,1,3-tri...	236	C18H20	003910-35-8	98
2	(Dimethylphenylsilyl)phenylacety...	236	C16H16Si	1000434-31-1	47
3	2-Pentene, 4-methyl-2,4-diphenyl-	236	C18H20	006258-73-7	41
4	4-Hydroxy-6-(trifluoromethyl)pyr...	236	C8H11F3N2OSi	1000476-16-0	38
5	Hexobarbital	236	C12H16N2O3	000056-29-1	38





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Instrument :
MSVOA_X
ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

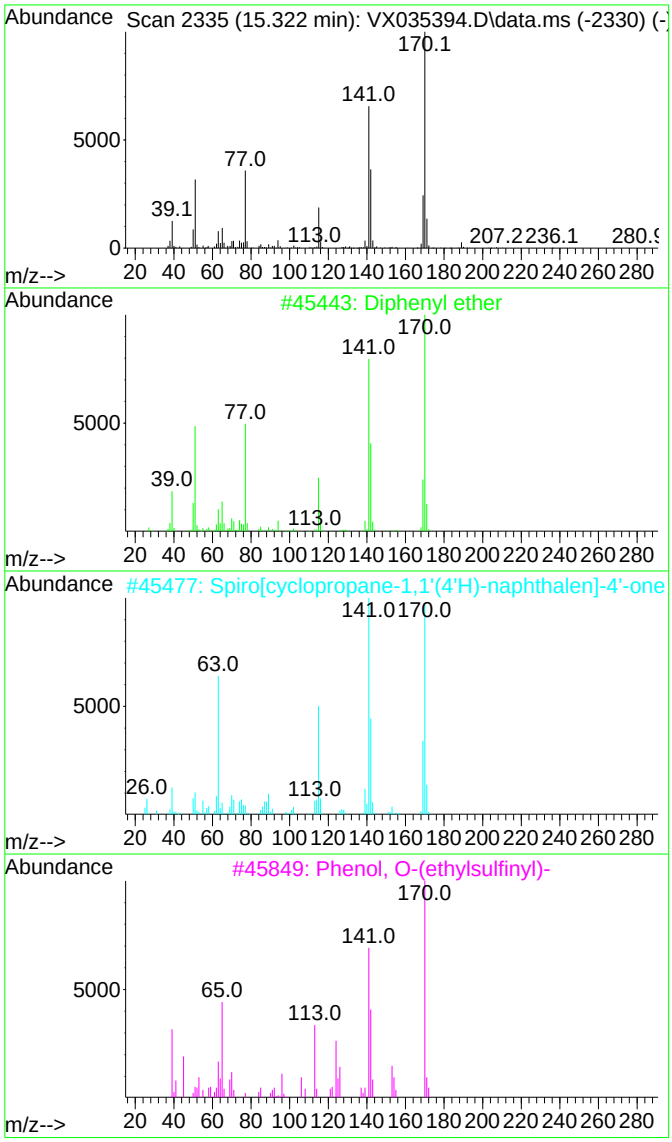
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

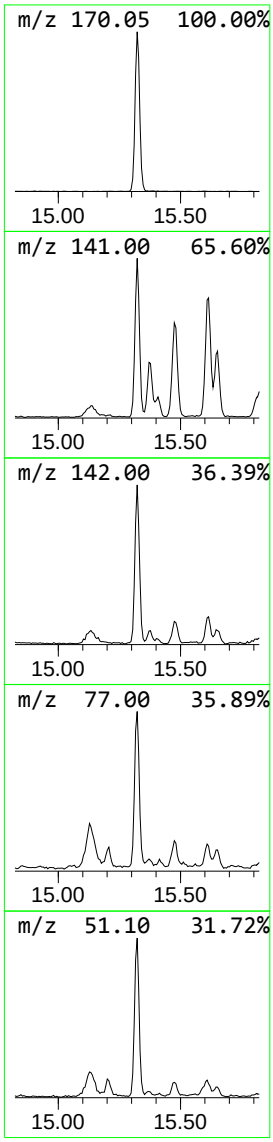
****

Peak Number 10 Diphenyl ether Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.323	12.47 ug/l	515297	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diphenyl ether	170	C12H10O	000101-84-8	96
2	Spiro[cyclopropane-1,1'(4'H)-nap...	170	C12H10O	033498-24-7	74
3	Phenol, O-(ethylsulfinyl)-	170	C8H10O2S	029634-40-0	50
4	2-Carbonitrile-8-hydroxyquinolin...	254	C15H14N2O2	1000463-12-5	43
5	Pyrazine, 2-(4-methylphenyl)-	170	C11H10N2	1000380-54-6	43





Tentatively Identified Compound (LSC) summary

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MSVOA_X
ClientSampleId :
SPA-4WC-23-04

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.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
Cyclohexane, pr...	10.854	10.3	ug/l	282537	3	10.055	1377250	50.0
Benzene, 1-ethy...	11.378	25.4	ug/l	1051480	4	12.024	2066870	50.0
Benzene, 1-ethy...	11.610	11.0	ug/l	453928	4	12.024	2066870	50.0
Undecane	12.421	13.2	ug/l	544333	4	12.024	2066870	50.0
Benzene, (1,1-d...	12.701	10.5	ug/l	432208	4	12.024	2066870	50.0
Benzene, 1,2,3,...	12.963	12.0	ug/l	496669	4	12.024	2066870	50.0
3-Phenylbut-1-ene	13.292	10.6	ug/l	436911	4	12.024	2066870	50.0
Benzocyclohepta...	14.634	13.5	ug/l	559448	4	12.024	2066870	50.0
1H-Indene, 2,3-...	15.134	32.4	ug/l	1337120	4	12.024	2066870	50.0
Diphenyl ether	15.323	12.5	ug/l	515297	4	12.024	2066870	50.0