LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0 % 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_U\Method\SFAMULM092022WMA.M
Title : VOC Analysis

Signal : TIC: VU050906.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.601	73	81	101	rVB2	1087905	1329618	7.39%	2.065%
2	1.906	168	176	195	rVB	183640	269308	1.50%	0.418%
3	2.514	363	365	368	rBV2	386	323	0.00%	0.001%
4	2.565	368	381	394	rBV	364911	600526	3.34%	0.933%
5	2.697	419	422	427	rBV3	435	419	0.00%	0.001%
6	2.787	443	450	456	rBV4	1307	2051	0.01%	0.003%
7	3.041	523	529	539	rVV4	1610	2623	0.01%	0.004%
8	3.179	560	572	584	rBV5	1241	2398	0.01%	0.004%
9	3.353	611	626	650	rVV2	143983	284199	1.58%	0.441%
10	3.581	695	697	702	rVB2	396	361	0.00%	0.001%
11	3.867	769	786	804	rBV	19464	46985	0.26%	0.073%
12	4.096	851	857	862	rVB2	353	465	0.00%	0.001%
13	4.154	867	875	879	rBV2	377	483	0.00%	0.001%
14	4.221	891	896	900	rVB3	442	395	0.00%	0.001%
15	4.298	914	920	928	rVB2	323	404	0.00%	0.001%
16	4.539	992	995	999	rVB2	637	477	0.00%	0.001%
17	4.665	1005	1034	1061	rBV2	2245622	5462450	30.37%	8.483%
18	5.060	1142	1157	1183	rVB	327183	713696	3.97%	1.108%
19	5.276	1222	1224	1226	rBV	674	371	0.00%	0.001%
20	5.379	1250	1256	1259	rBV5	2256	2725	0.02%	0.004%
21	5.453	1275	1279	1283	rBV2	639	546	0.00%	0.001%
22	5.527	1299	1302	1310	rVB2	403	380	0.00%	0.001%
23	5.723	1338	1363	1382	rBV2	668152	1795784	9.99%	2.789%
24	5.980	1432	1443	1444	rBV5	1736	2159	0.01%	0.003%
25	6.044	1458	1463	1466	rBV4	728	814	0.00%	0.001%
26	6.115	1481	1485	1489	rBV4	588	620	0.00%	0.001%
27	6.157	1494	1498	1505	rVB4	587	883	0.00%	0.001%
28	6.247	1512	1526	1548	rBV	509698	965512	5.37%	1.499%
29	6.453	1588	1590	1596	rVB3	718	581	0.00%	0.001%
30	6.533	1601	1615	1627	rVB10	9747	23487	0.13%	0.036%
31	6.690	1646	1664	1678	rBV	430703	834554	4.64%	1.296%
32	6.912	1730	1733	1738	rVB3	544	365	0.00%	0.001%
33	6.941	1738	1742	1745	rVB2	606	487	0.00%	0.001%
34	6.964	1745	1749	1750	rBV2	530	407	0.00%	0.001%
35	6.980	1750	1754	1762	rVB6	1308	2005	0.01%	0.003%
36	7.054	1773	1777	1779	rBV2	659	485	0.00%	0.001%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0 % 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_U\Method\SFAMULM092022WMA.M
Title : VOC Analysis

37	7.070	1779	1782	1790	rVB5	1056	1099	0.01%	0.002%
38	7.144	1800	1805	1807	rBV	746	557	0.00%	0.001%
39	7.166	1807	1812	1813	rBV2	1266	1009	0.01%	0.002%
40	7.234	1828	1833	1835	rBV5	1132	1094	0.01%	0.002%
41	7.379	1871	1878	1880	rBV2	324	346	0.00%	0.001%
42	7.565	1921	1936	1953	rBV	211505	375626	2.09%	0.583%
43	7.793	2003	2007	2013	rVB3	516	470	0.00%	0.001%
44	7.896	2016	2039	2054	rBV	777285	1353505	7.53%	2.102%
45	8.031	2073	2081	2094	rVB8	7618	15209	0.08%	0.024%
46	8.096	2098	2101	2110	rVB5	4935	5490	0.03%	0.009%
47	8.176	2113	2126	2138	rBV2	153829	258970	1.44%	0.402%
48	8.237	2139	2145	2158	rVB4	21765	40417	0.22%	0.063%
49	8.369	2176	2186	2198	rVB4	8598	15254	0.08%	0.024%
50	8.417	2198	2201	2203	rBV3	624	499	0.00%	0.001%
51	8.475	2209	2219	2231	rBV6	3949	7373	0.04%	0.011%
52	8.536	2233	2238	2241	rBV5	1733	2211	0.01%	0.003%
53	8.633	2253	2268	2301	rBV2	576652	1126515	6.26%	1.749%
54	8.803	2311	2321	2330	rBV2	13169	26896	0.15%	0.042%
55	8.864	2331	2340	2349	rVB2	44584	69161	0.38%	0.107%
56	8.922	2349	2358	2368	rBV2	46867	79820	0.44%	0.124%
57	9.070	2391	2404	2411	rBV6	9837	17282	0.10%	0.027%
58	9.115	2411	2418	2424	rVV5	5972	8520	0.05%	0.013%
59	9.163	2425	2433	2442	rVV6	24838	41569	0.23%	0.065%
60	9.208	2443	2447	2457	rVB3	11762	17079	0.09%	0.027%
61	9.295	2457	2474	2478	rBV9	4742	9389	0.05%	0.015%
62	9.333	2479	2486	2498	rVB4	12064	19951	0.11%	0.031%
63	9.417	2499	2512	2528	rBV	727559	1204310	6.70%	1.870%
64	9.510	2534	2541	2548	rVB2	32691	46155	0.26%	0.072%
65	9.568	2549	2559	2572	rVB	177029	286219	1.59%	0.444%
66	9.697	2573	2599	2612	rBV3	207270	466226	2.59%	0.724%
67	9.883	2644	2657	2668	rBV6	33280	74409	0.41%	0.116%
68	9.951	2674	2678	2685	rVB5	4314	5136	0.03%	0.008%
69	10.034	2689	2704	2713	rBV6	73189	208405	1.16%	0.324%
70	10.269	2758	2777	2786	rBV7	137679	387279	2.15%	0.601%
71	10.356	2797	2804	2809	rBV4	77164	129207	0.72%	0.201%
72	10.481	2834	2843	2852	rVB	337154	520351	2.89%	0.808%
73	10.697	2904	2910	2917	rBV3	57524	73822	0.41%	0.115%
74	10.755	2918	2928	2938	rBV	674120	1086812	6.04%	1.688%
75	10.809	2940	2945	2963	rVB7	183195	313165	1.74%	0.486%
76	10.906	2965	2975	2984	rBV	636507	988524	5.50%	1.535%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0 % 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_U\Method\SFAMULM092022WMA.M
Title : VOC Analysis

77	11.021	3001	3011	3020	rBV	1239038	1901051	10.57%	2.952%
78	11.292	3084	3095	3113	rBV	1664955	2715273	15.10%	4.217%
79	11.417	3127	3134	3139	rBV2	269213	371717	2.07%	0.577%
80	11.465	3139	3149	3173	rVB	11797592	17984659	100.00%	27.928%
81	11.809	3243	3256	3269	rVB3	847838	1772663	9.86%	2.753%
82	11.893	3271	3282	3298	rVB	2919083	4456908	24.78%	6.921%
83	12.005	3312	3317	3329	rVB	231246	338579	1.88%	0.526%
84	12.095	3335	3345	3357	rVB2	1013296	1569322	8.73%	2.437%
85	12.189	3362	3374	3392	rVB4	1179513	2390942	13.29%	3.713%
86	12.301	3401	3409	3421	rVB	253166	391259	2.18%	0.608%
87	12.375	3424	3432	3447	rVB	283624	495287	2.75%	0.769%
88	12.462	3449	3459	3464	rBV	752749	1159812	6.45%	1.801%
89	12.568	3484	3492	3500	rBV	923022	1311102	7.29%	2.036%
90	12.684	3517	3528	3545	rVB3	646929	1523950	8.47%	2.367%
91	12.870	3580	3586	3597	rVB3	246217	415587	2.31%	0.645%
92	12.970	3610	3617	3625	rVB	375170	533264	2.97%	0.828%
93	13.025	3625	3634	3650	rVB	790020	1250390	6.95%	1.942%
94	13.108	3653	3660	3678	rVB5	83573	146523	0.81%	0.228%
95	13.198	3681	3688	3696	rBV	41741	64110	0.36%	0.100%
96	13.401	3745	3751	3756	rBV4	36035	55084	0.31%	0.086%
97	13.449	3757	3766	3783	rVB2	659460	1126185	6.26%	1.749%
98	13.623	3811	3820	3836	rVB	258132	413546	2.30%	0.642%
99	14.086	3955	3964	3984	rVB	187871	331673	1.84%	0.515%
100	15.410	4368	4376	4388	rVB	25579	42700	0.24%	0.066%

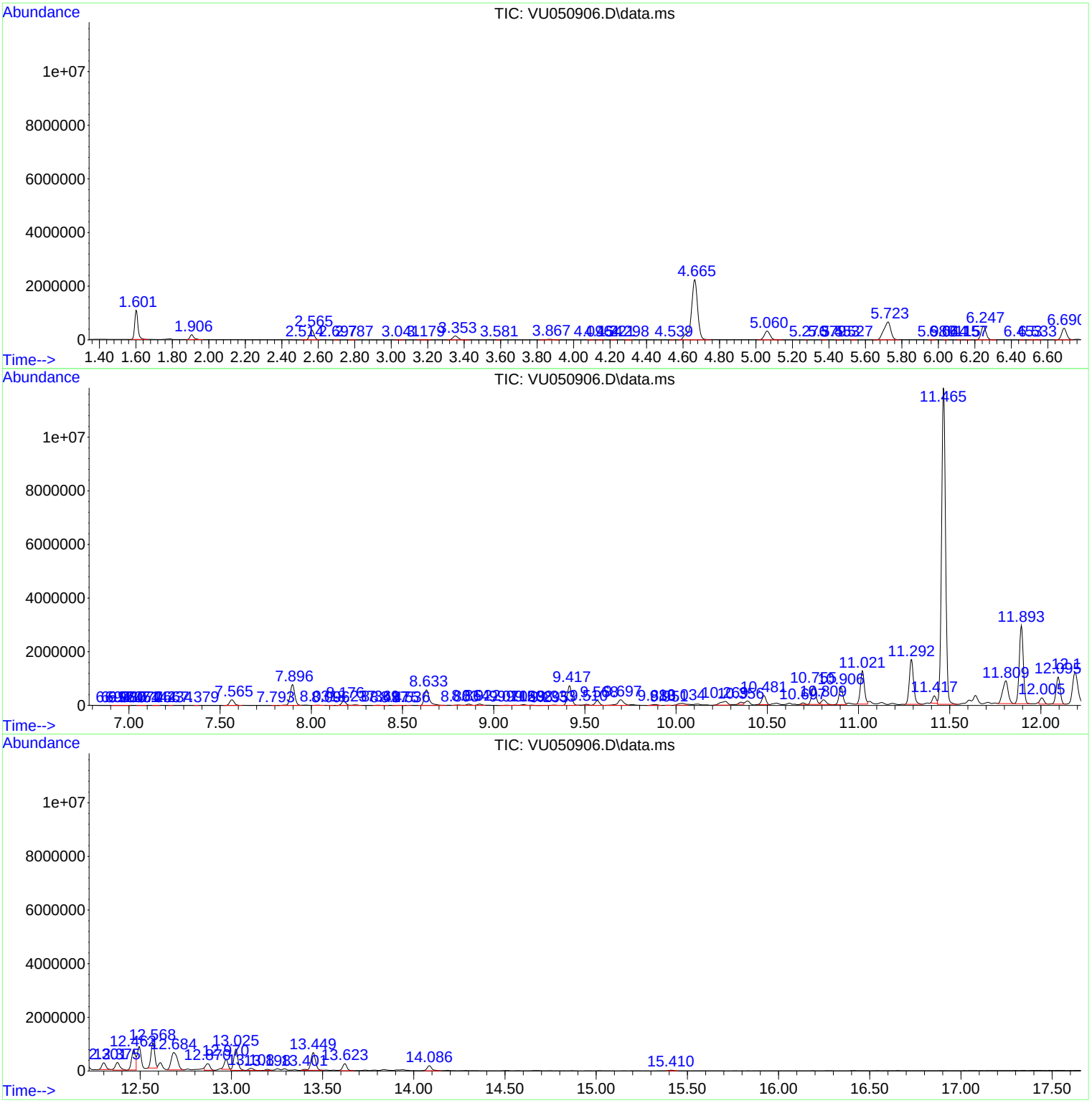
Sum of corrected areas: 64396263

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

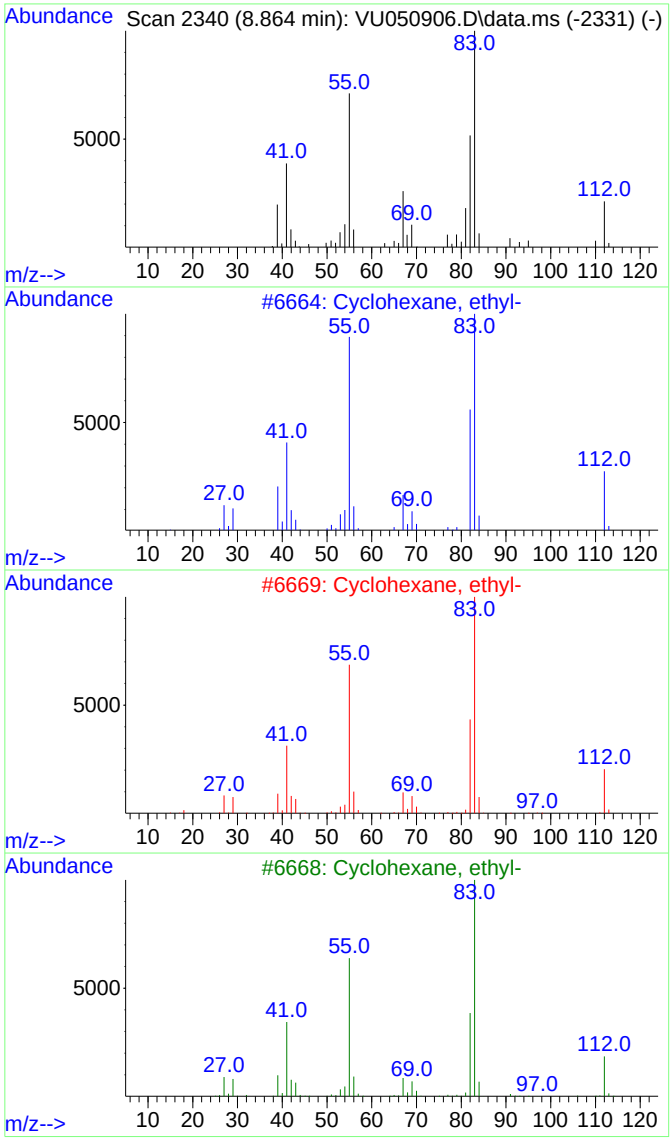
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

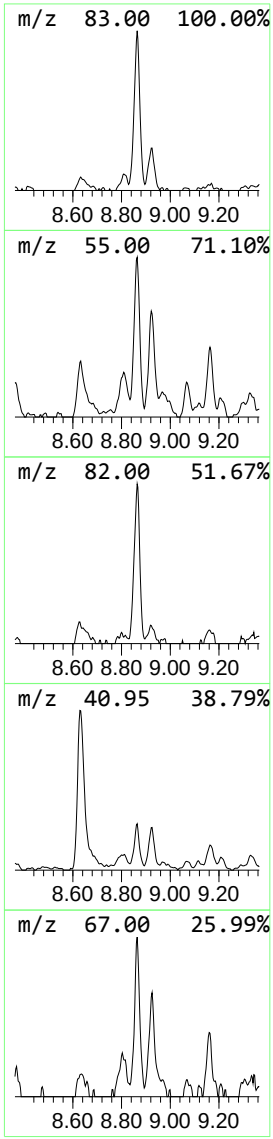
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 (DEL) Alkane: Cyclic8.864 Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.864	2.87 ug/L	69161	Chlorobenzene-d5	9.417

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, ethyl-	112	C8H16	001678-91-7	93
2	Cyclohexane, ethyl-	112	C8H16	001678-91-7	83
3	Cyclohexane, ethyl-	112	C8H16	001678-91-7	83
4	Cyclohexane, ethyl-	112	C8H16	001678-91-7	64
5	Cyclohexane, ethyl-	112	C8H16	001678-91-7	53





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

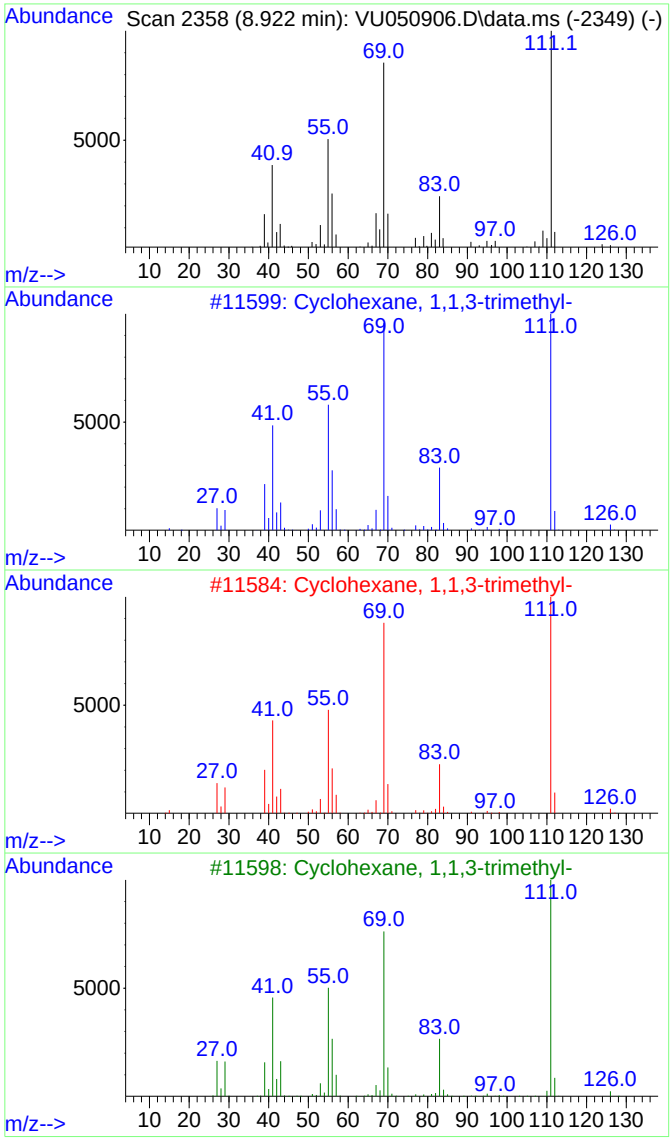
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

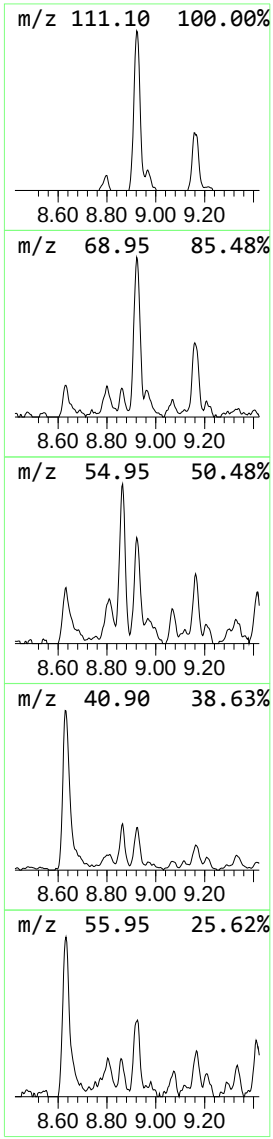
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 (DEL) Alkane: Cyclic8.922 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.922	3.31 ug/L	79820	Chlorobenzene-d5	9.417

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	93
2	Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	87
3	Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	74
4	1,1,4-Trimethylcyclohexane	126	C9H18	007094-27-1	64
5	Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-26-2	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

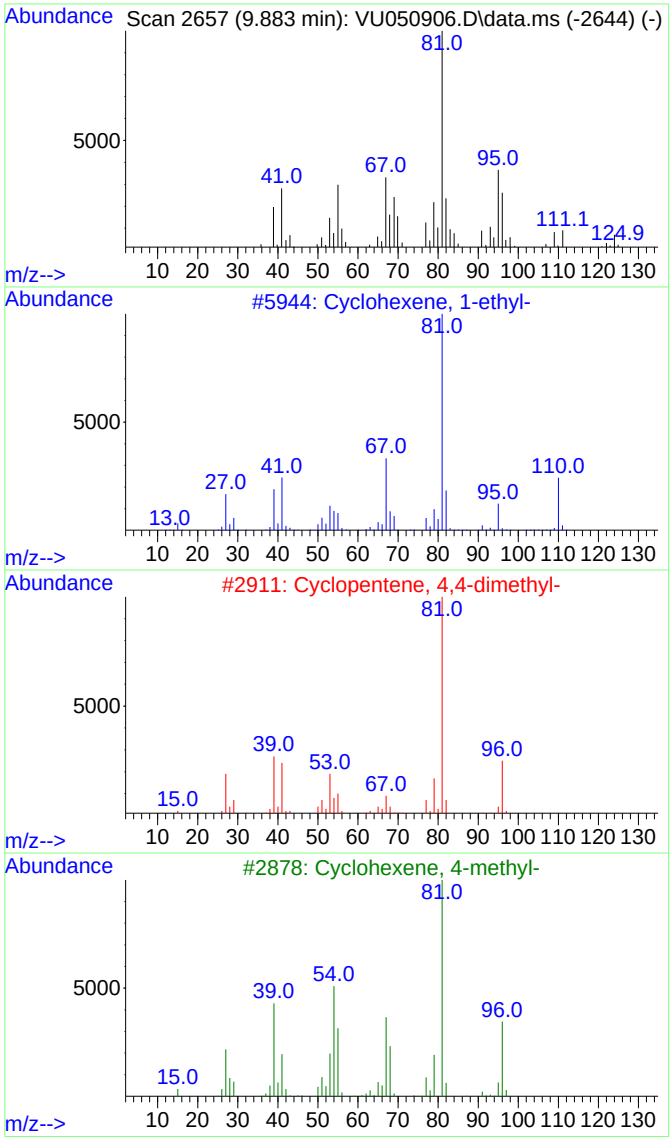
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

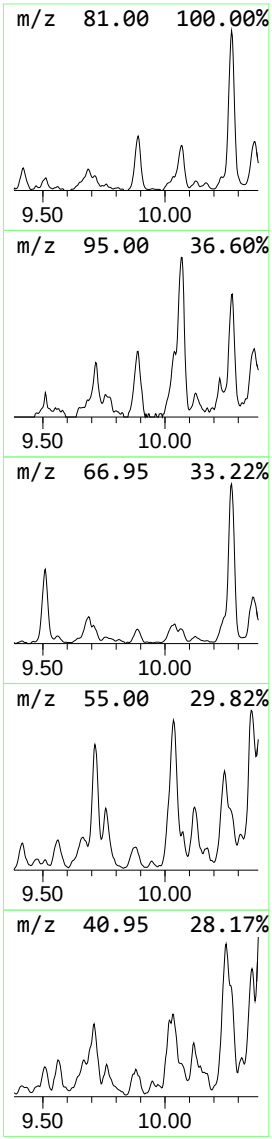
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Cyclohexene, 1-ethyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.883	3.09 ug/L	74409	Chlorobenzene-d5	9.417

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexene, 1-ethyl-	110	C8H14	001453-24-3	53
2	Cyclopentene, 4,4-dimethyl-	96	C7H12	019037-72-0	52
3	Cyclohexene, 4-methyl-	96	C7H12	000591-47-9	52
4	1-Ethyl-5-methylcyclopentene	110	C8H14	097797-57-4	50
5	Cyclobutane, (1-methylethylidene)-	96	C7H12	001528-22-9	49





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

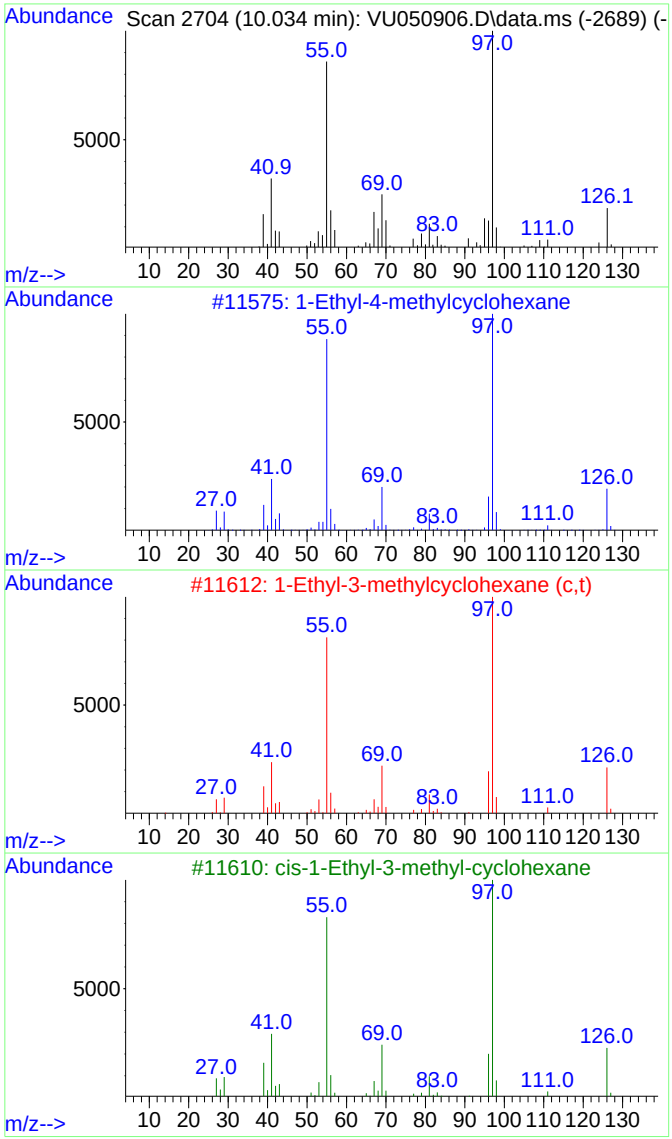
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

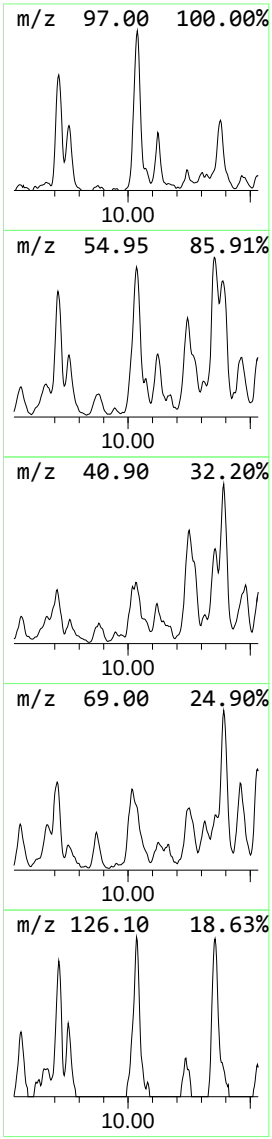
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 (DEL) Alkane: Cyclic10.034 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.034	8.65 ug/L	208405	Chlorobenzene-d5	9.417

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	83
2	1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	81
3	cis-1-Ethyl-3-methyl-cyclohexane	126	C9H18	019489-10-2	74
4	Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	72
5	Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	72





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

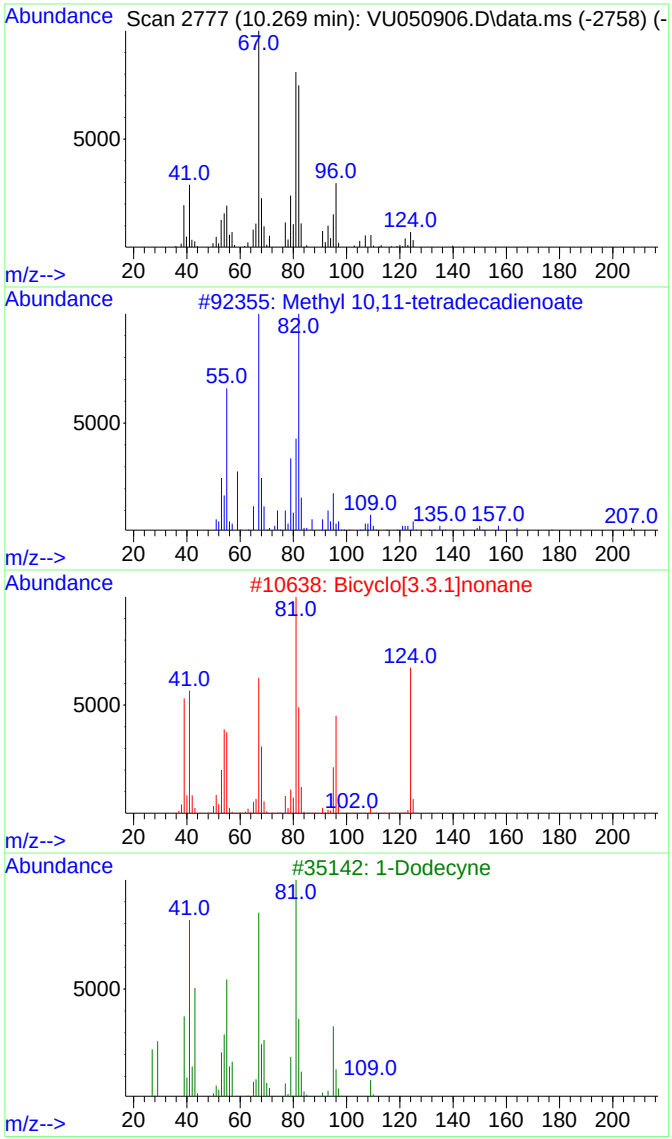
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

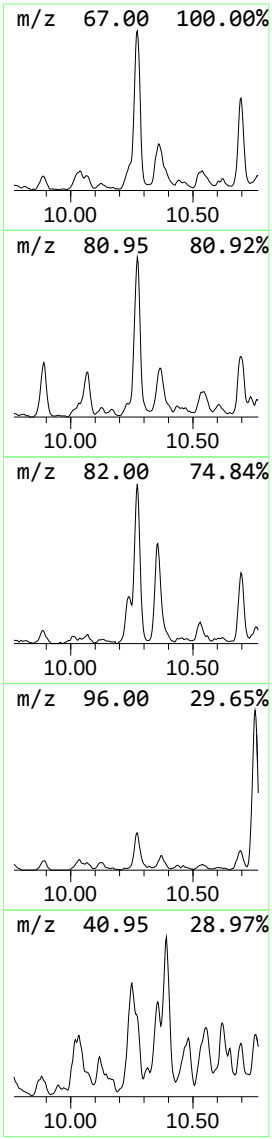
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Methyl 10,11-tetradecadienoate Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.269	16.08 ug/L	387279	Chlorobenzene-d5	9.417

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl 10,11-tetradecadienoate	238	C15H26O2	1000336-31-8	50
2	Bicyclo[3.3.1]nonane	124	C9H16	000280-65-9	46
3	1-Dodecyne	166	C12H22	000765-03-7	43
4	C2H5CH=CHCH=CH2	82	C6H10	000592-48-3	43
5	Cyclohexene, 3-methyl-	96	C7H12	000591-48-0	43





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

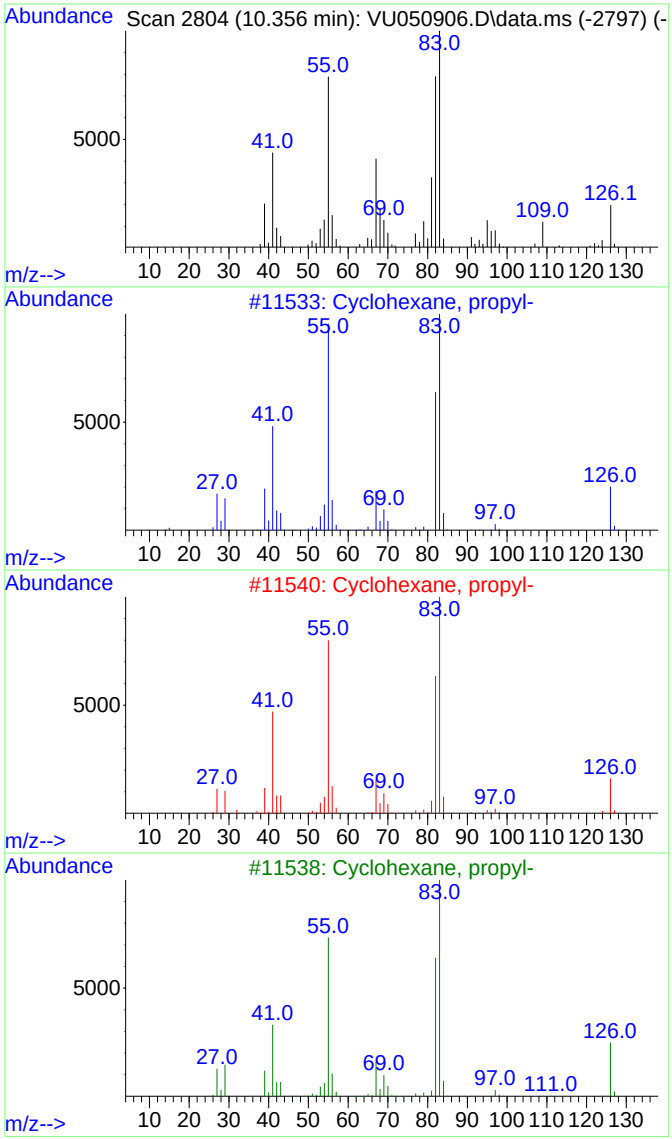
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

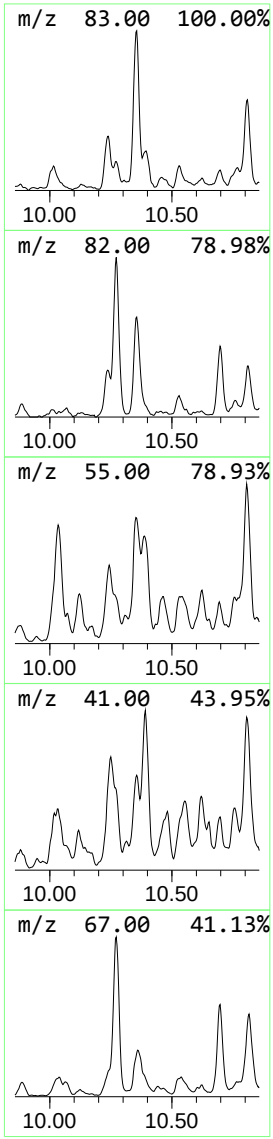
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 (DEL) Alkane: Cyclic10.356 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.356	5.36 ug/L	129207	Chlorobenzene-d5	9.417

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, propyl-	126	C9H18	001678-92-8	76
2	Cyclohexane, propyl-	126	C9H18	001678-92-8	72
3	Cyclohexane, propyl-	126	C9H18	001678-92-8	64
4	Cyclohexane, octyl-	196	C14H28	001795-15-9	53
5	Vinylcyclohexyl ether	126	C8H14O	002182-55-0	53





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

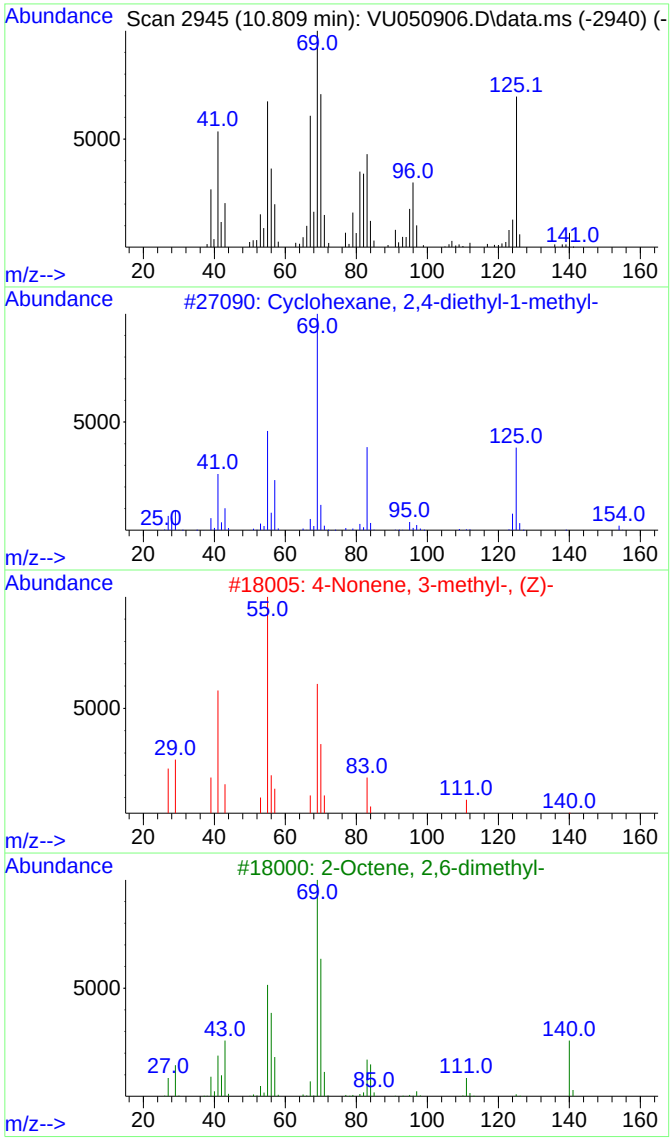
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

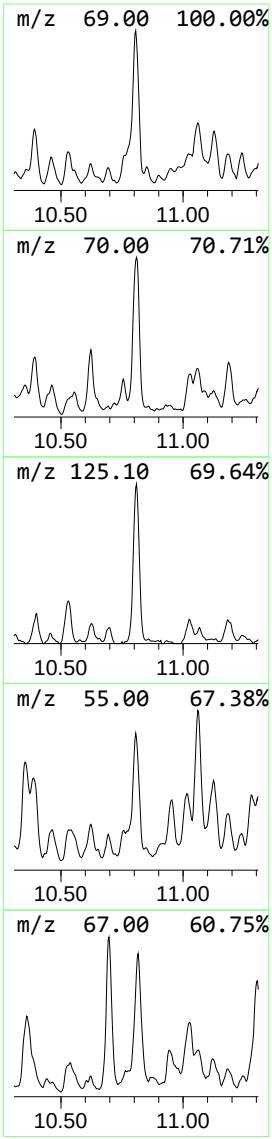
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Cyclic10.809 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.809	8.83 ug/L	313165	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 2,4-diethyl-1-methyl-	154	C11H22	061142-70-9	50
2	4-Nonene, 3-methyl-, (Z)-	140	C10H20	063830-69-3	43
3	2-Octene, 2,6-dimethyl-	140	C10H20	004057-42-5	41
4	4N-Methylcytosine	125	C5H7N3O	006220-47-9	41
5	1-Hexene, 3,3-dimethyl-	112	C8H16	003404-77-1	35





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

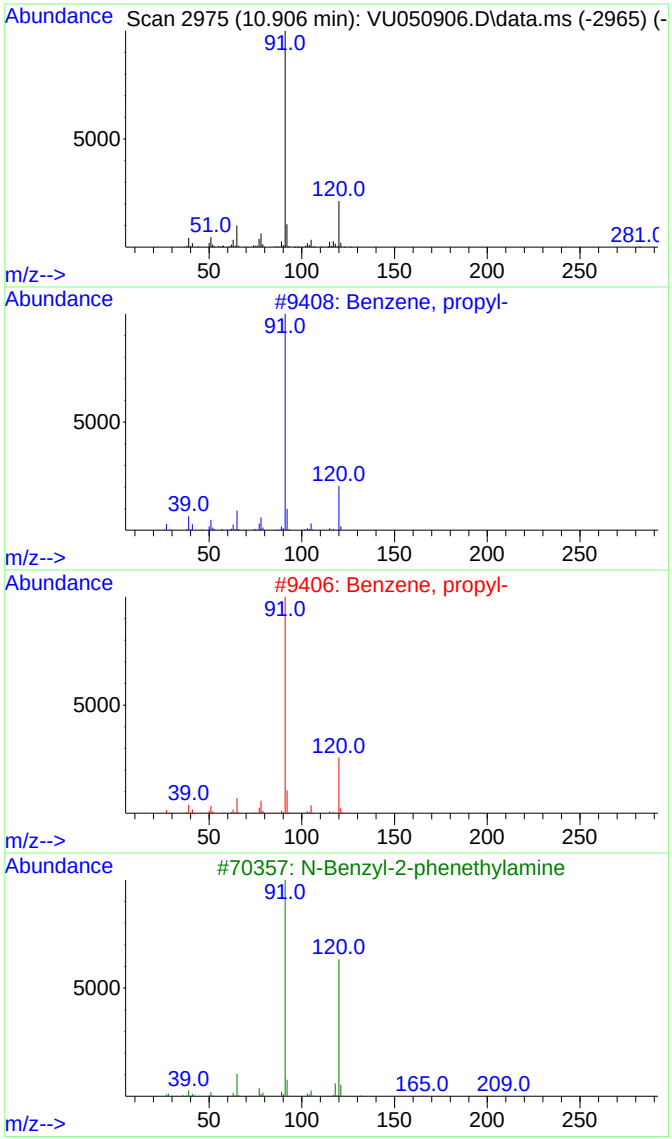
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

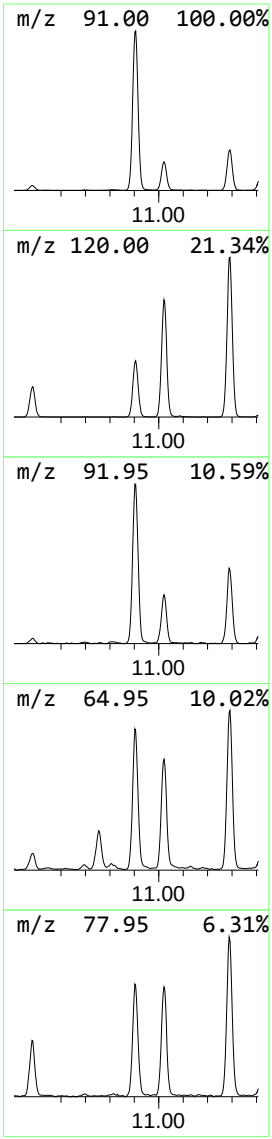
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, propyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.906	27.88 ug/L	988524	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	90
2	Benzene, propyl-	120	C9H12	000103-65-1	87
3	N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	78
4	Benzene, propyl-	120	C9H12	000103-65-1	78
5	Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	72





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

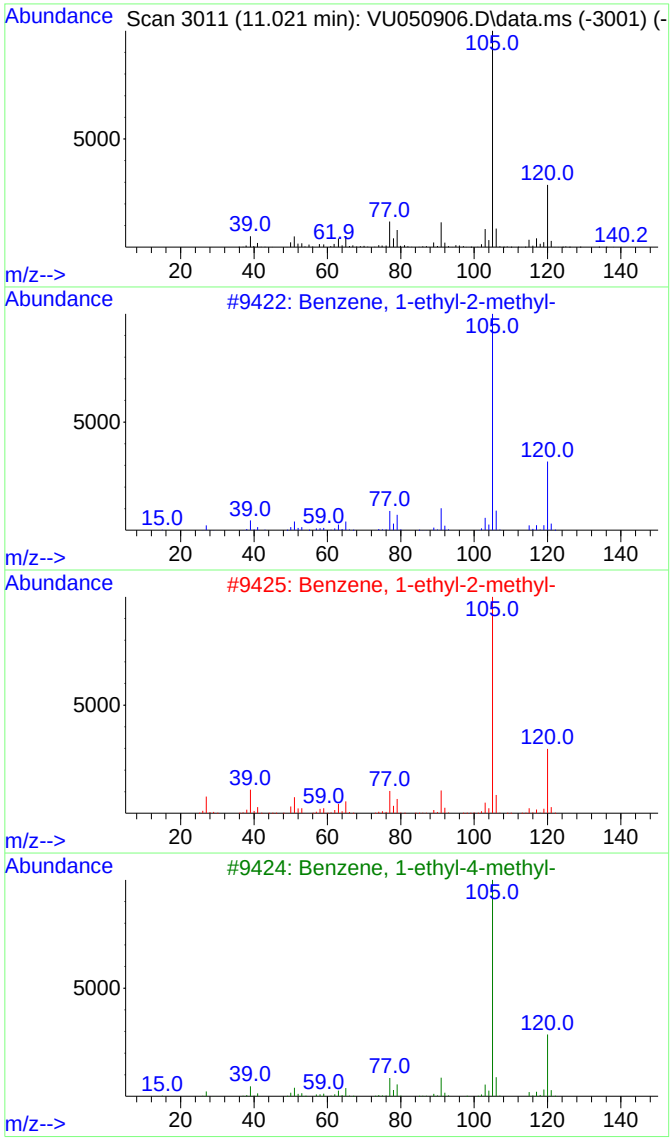
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

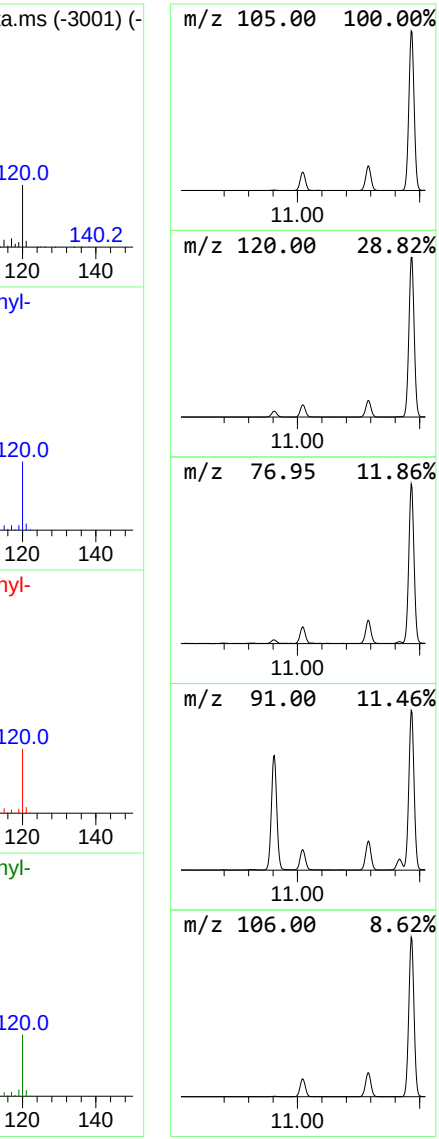
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.022	53.62 ug/L	1901050	1,4-Dichlorobenzene-d4	11.809

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-2-methyl-	120	C9H12	000611-14-3	93
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
4	Mesitylene	120	C9H12	000108-67-8	90
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

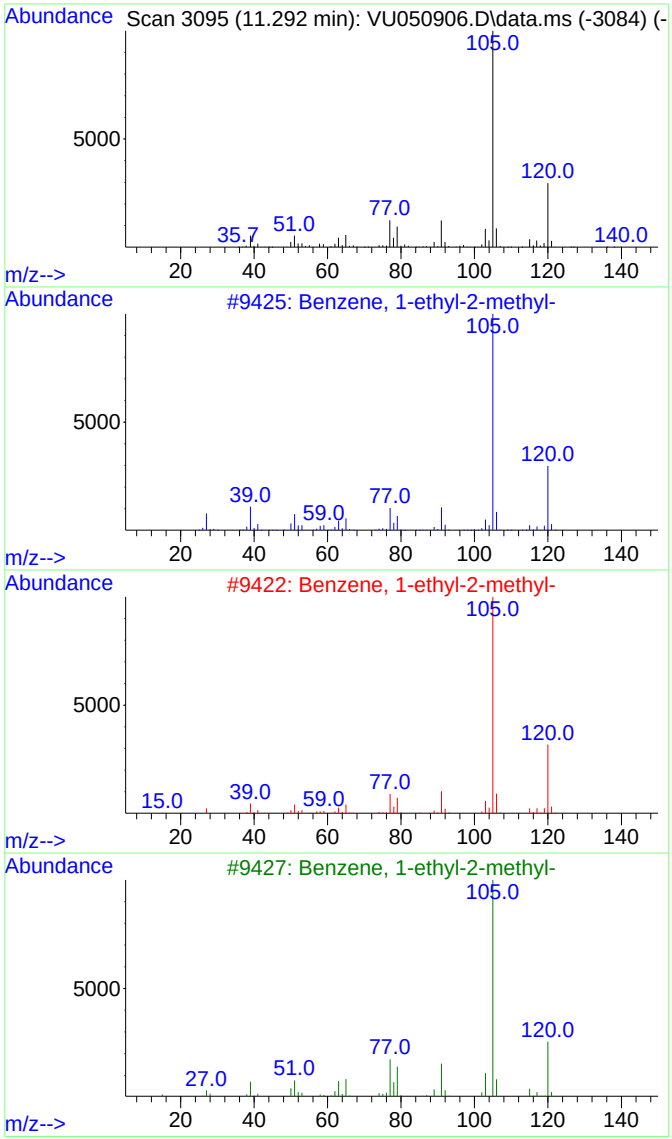
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

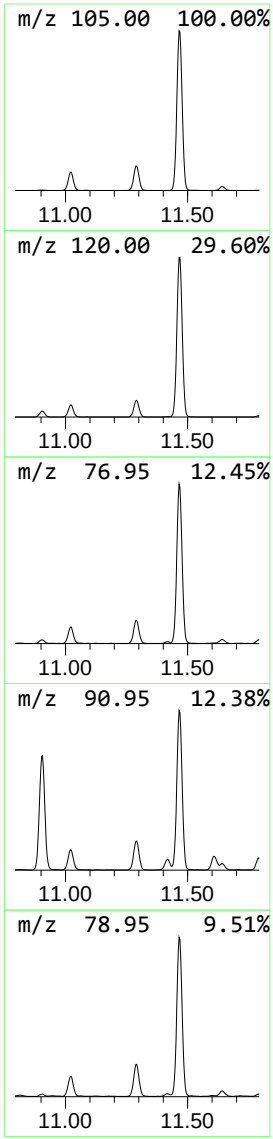
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Benzene, 1-ethyl-3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.292	76.59 ug/L	2715270	1,4-Dichlorobenzene-d4	11.809

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-2-methyl-	120	C9H12	000611-14-3	94
3	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

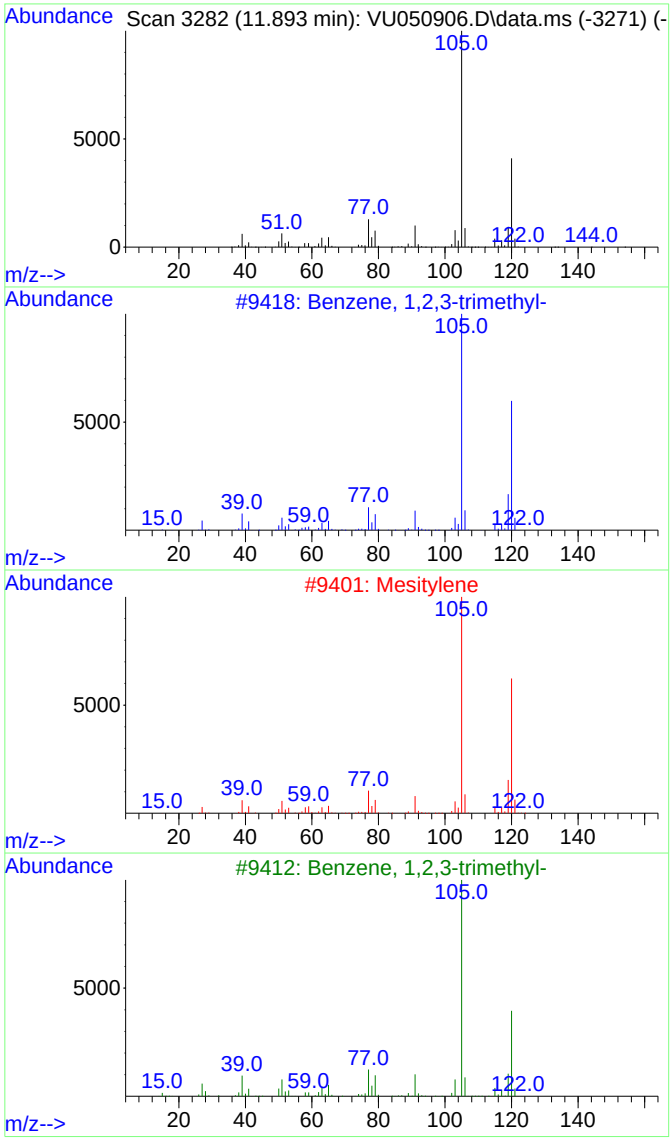
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

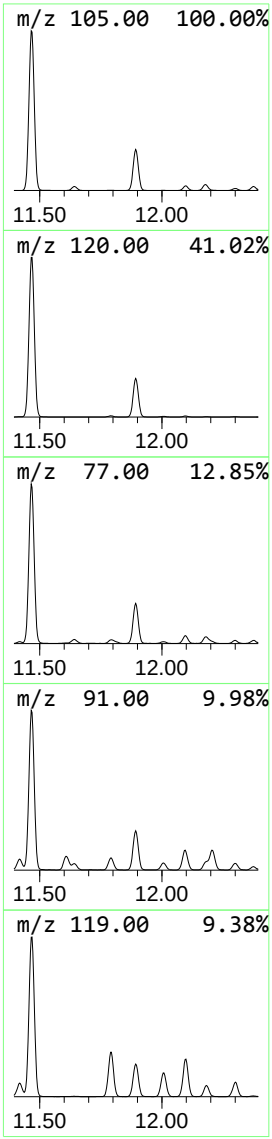
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Benzene, 1,2,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.893	125.71 ug/L	4456910	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2	Mesitylene	120	C9H12	000108-67-8	94
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

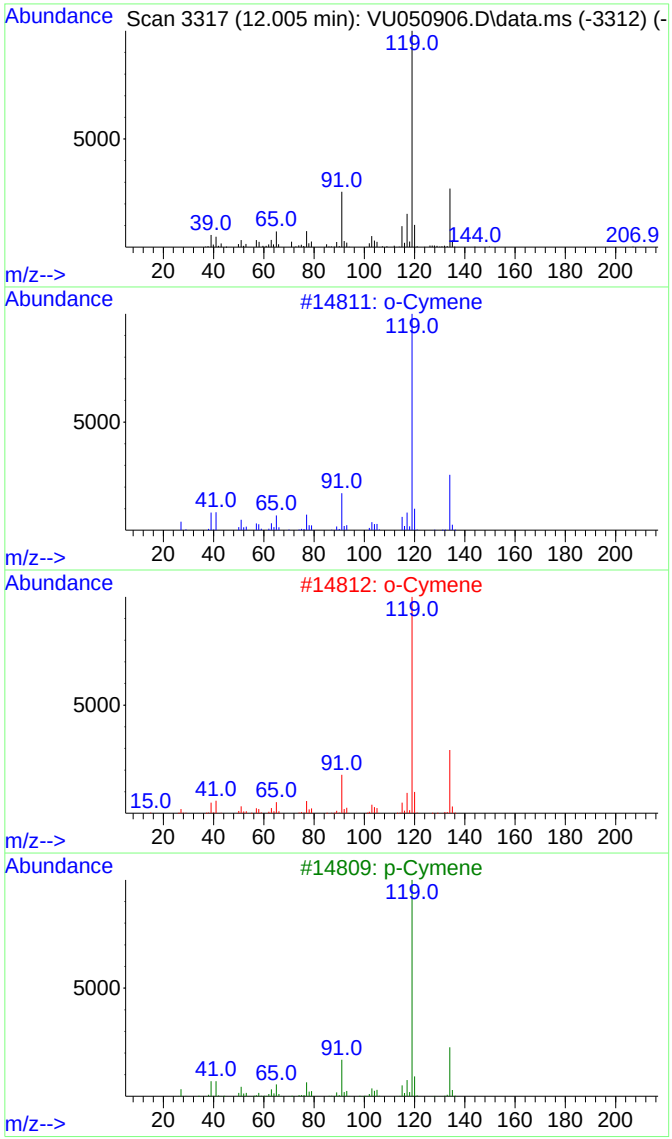
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

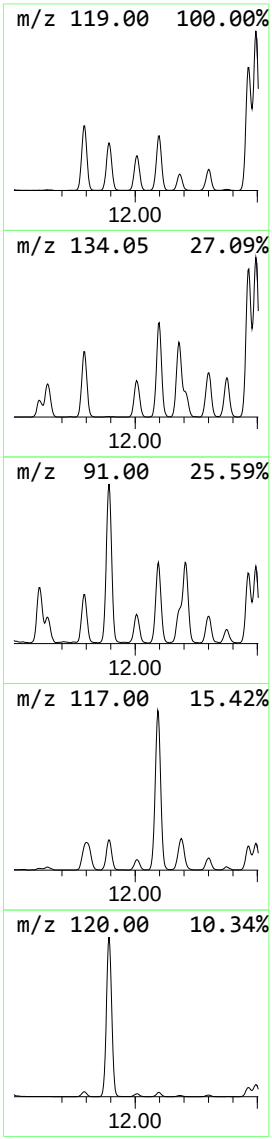
****

Peak Number 13 p-Cymene Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.005	9.55 ug/L	338579	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	97
2	o-Cymene	134	C10H14	000527-84-4	95
3	p-Cymene	134	C10H14	000099-87-6	94
4	o-Cymene	134	C10H14	000527-84-4	94
5	o-Cymene	134	C10H14	000527-84-4	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

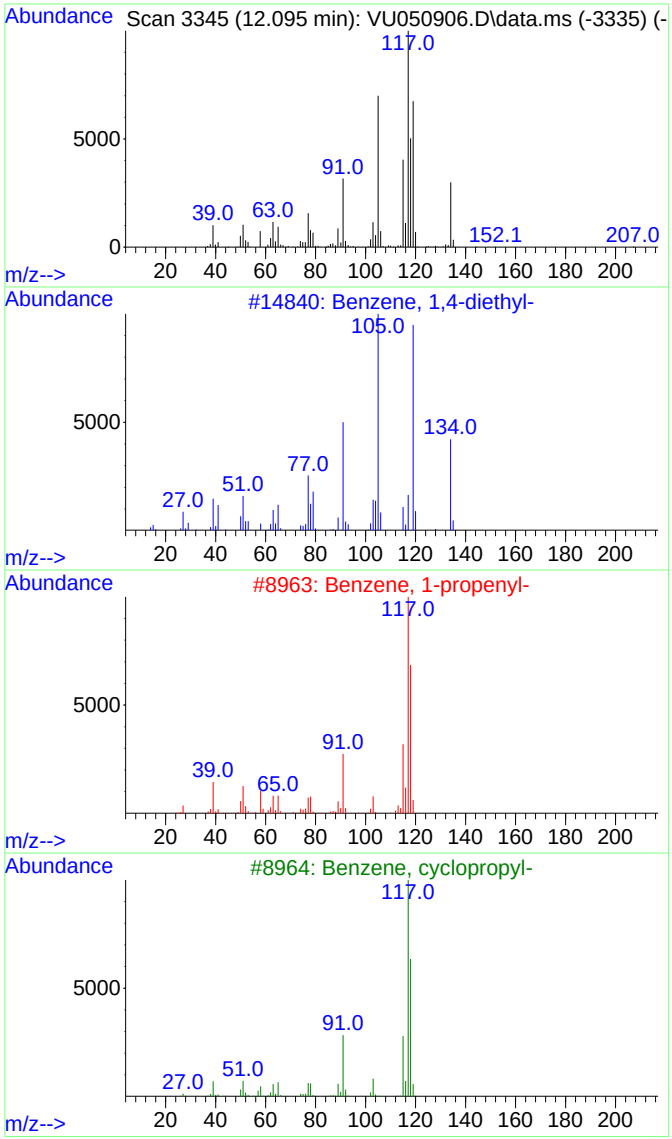
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

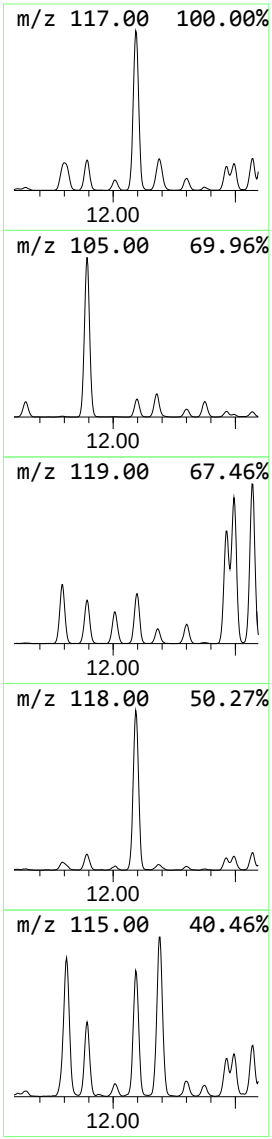
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Benzene, 1,4-diethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.095	44.26 ug/L	1569320	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	55
2	Benzene, 1-propenyl-	118	C9H10	000637-50-3	55
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	55
4	Benzene, 2-propenyl-	118	C9H10	000300-57-2	55
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	55





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

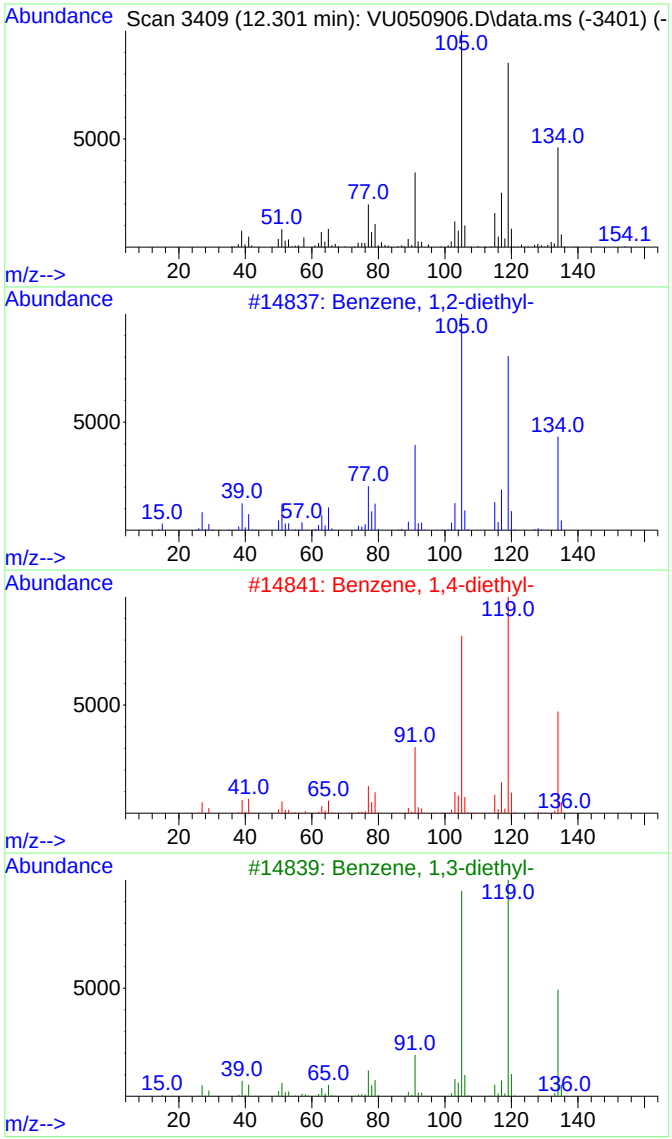
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

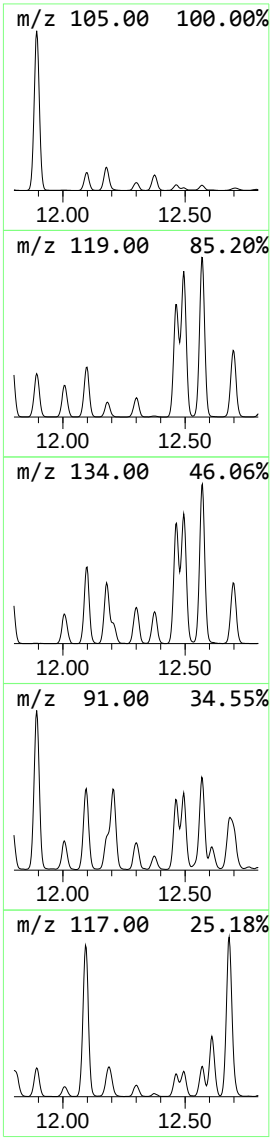
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Benzene, 1,2-diethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.301	11.04 ug/L	391259	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
2	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	93
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	90
4	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	90
5	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

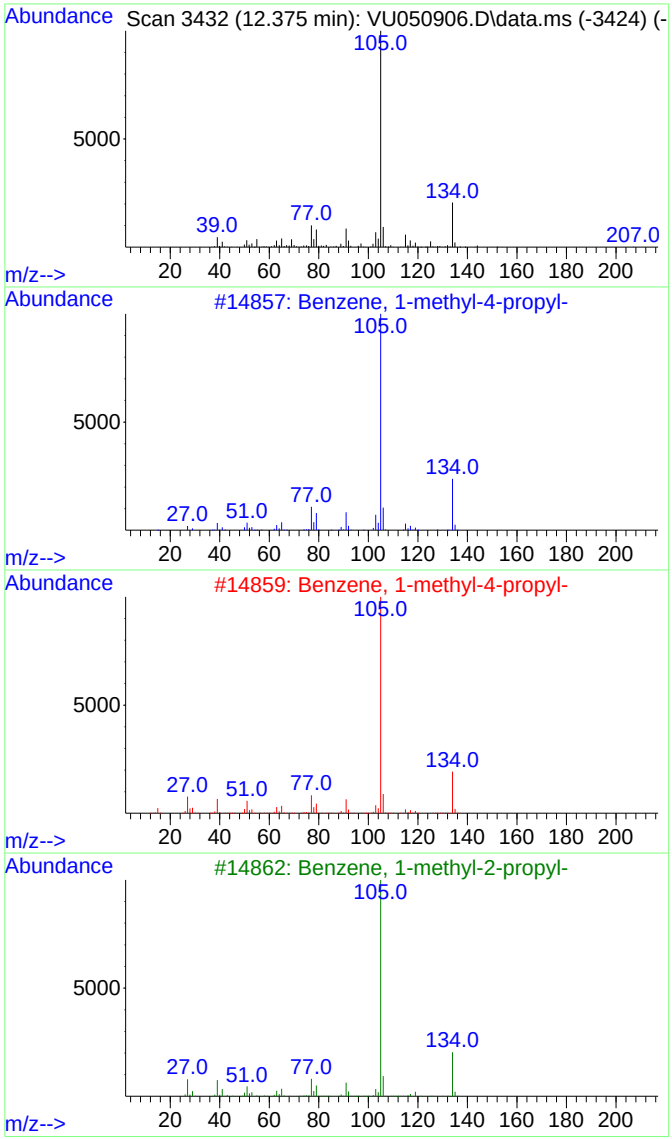
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

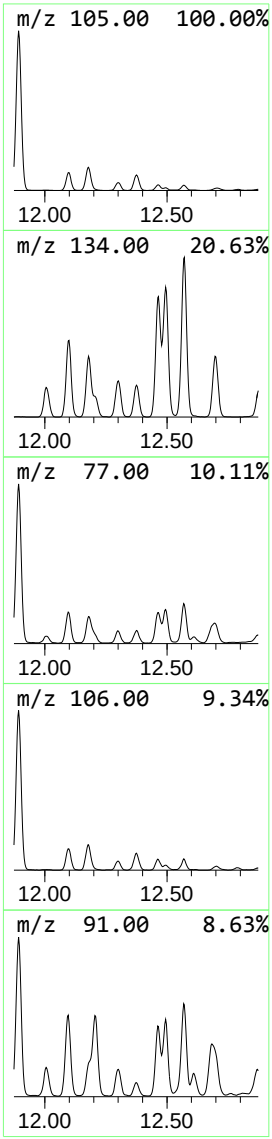
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Benzene, 1-methyl-4-propyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.375	13.97 ug/L	495287	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	95
2	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
3	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91
4	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91
5	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

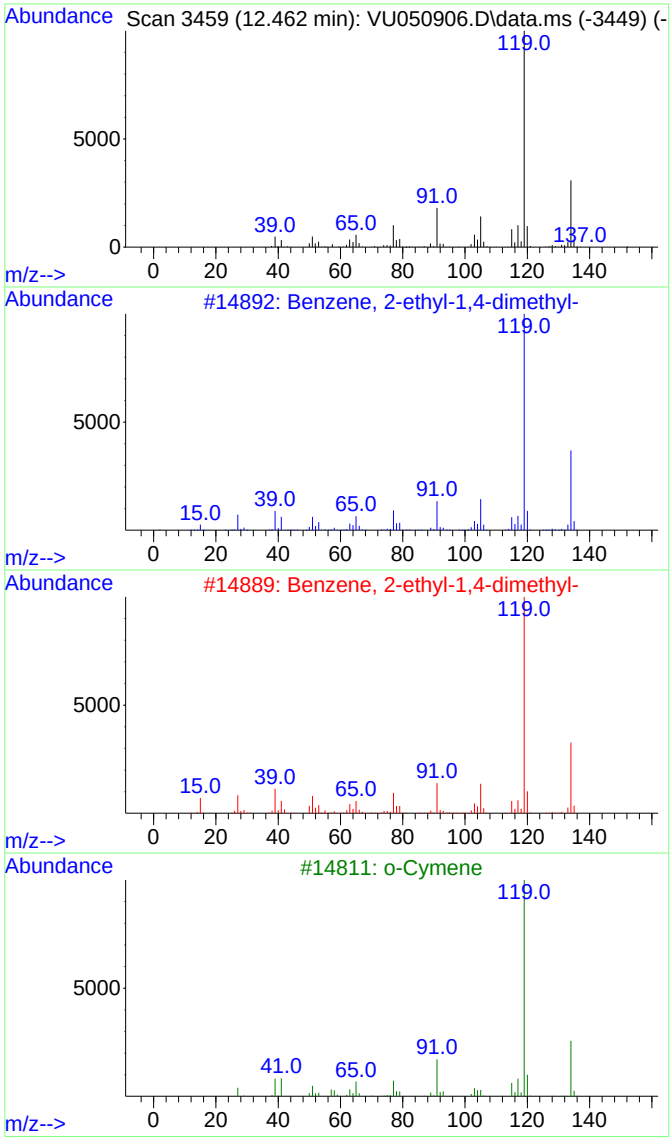
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

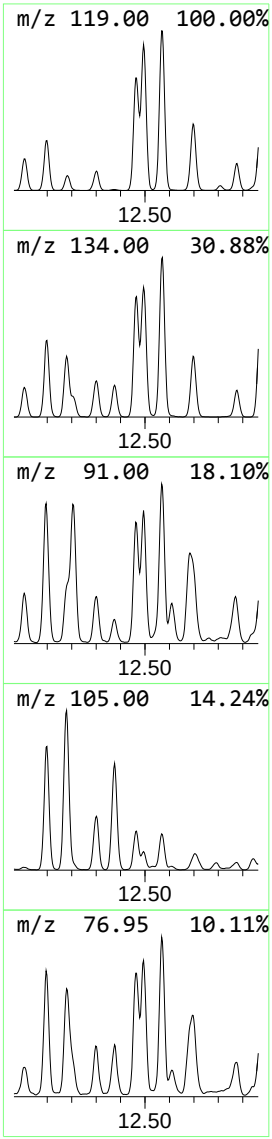
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.462	32.71 ug/L	1159810	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3	o-Cymene	134	C10H14	000527-84-4	95
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

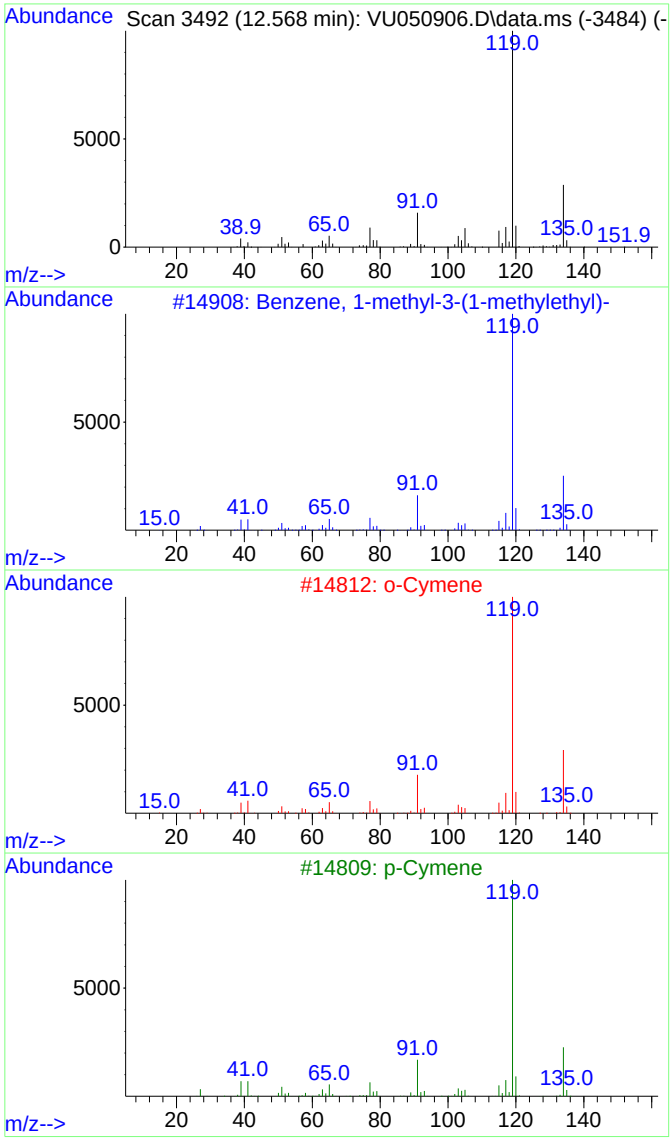
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

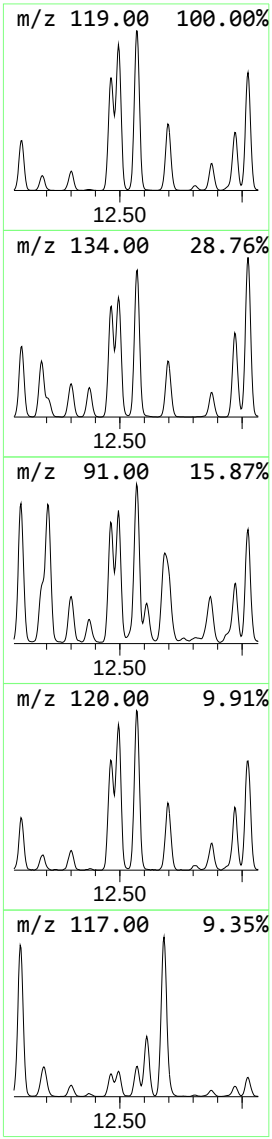
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Benzene, 1-methyl-3-(1-meth... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.568	36.98 ug/L	1311100	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
2	o-Cymene	134	C10H14	000527-84-4	95
3	p-Cymene	134	C10H14	000099-87-6	95
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

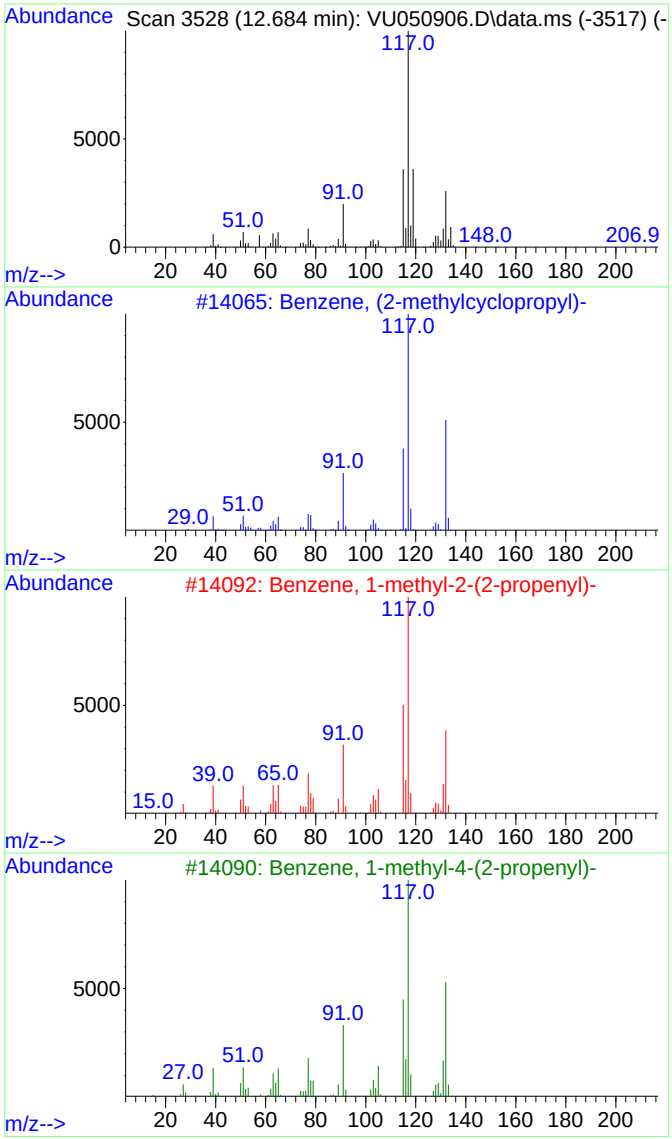
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

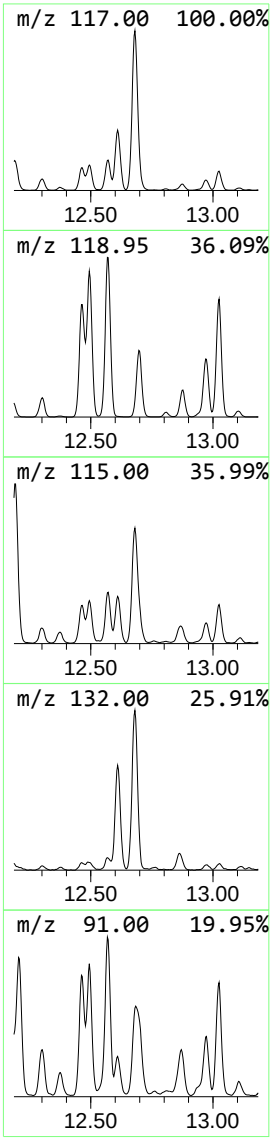
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Benzene, (2-methylcycloprop... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.684	42.98 ug/L	1523950	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	76
2	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	74
3	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	74
4	Benzene, 2-butenyl-	132	C10H12	001560-06-1	72
5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	72





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

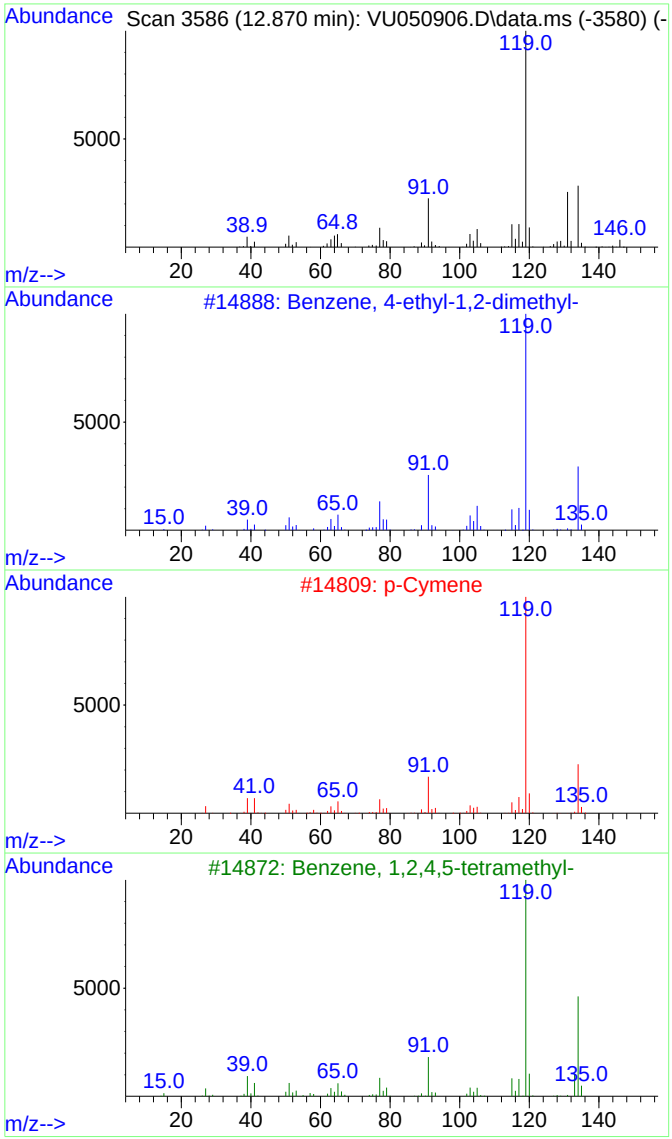
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

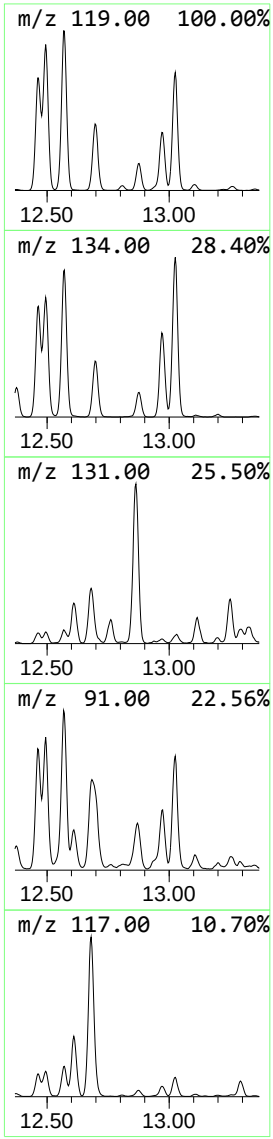
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.870	11.72 ug/L	415587	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
2	p-Cymene	134	C10H14	000099-87-6	93
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	93
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	81
5	o-Cymene	134	C10H14	000527-84-4	81





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

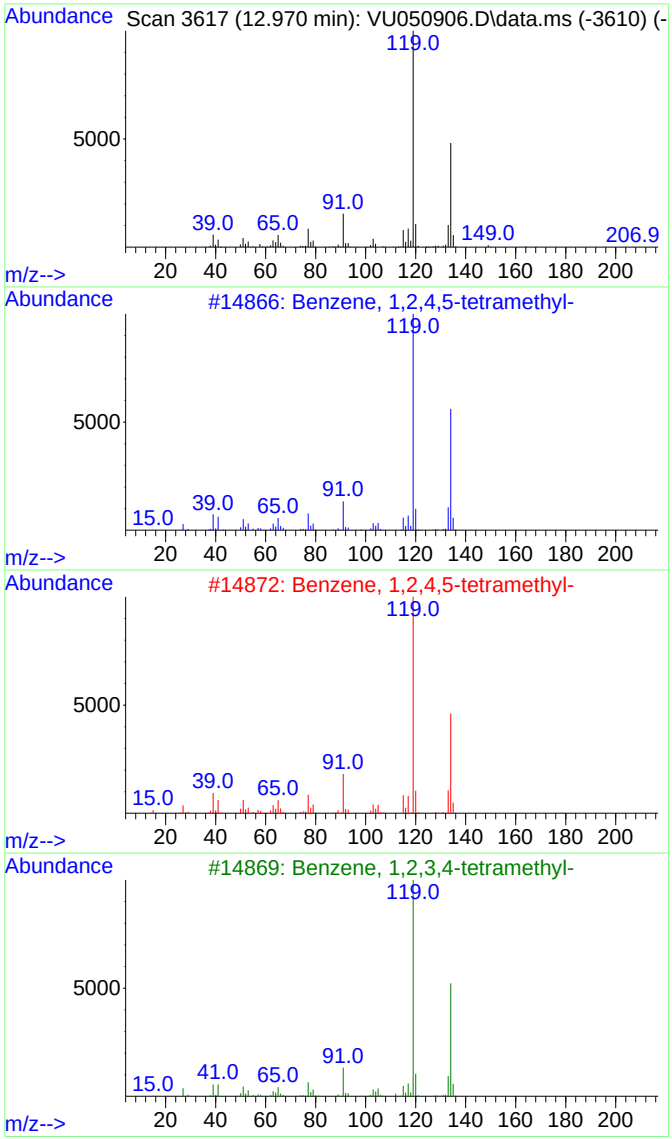
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

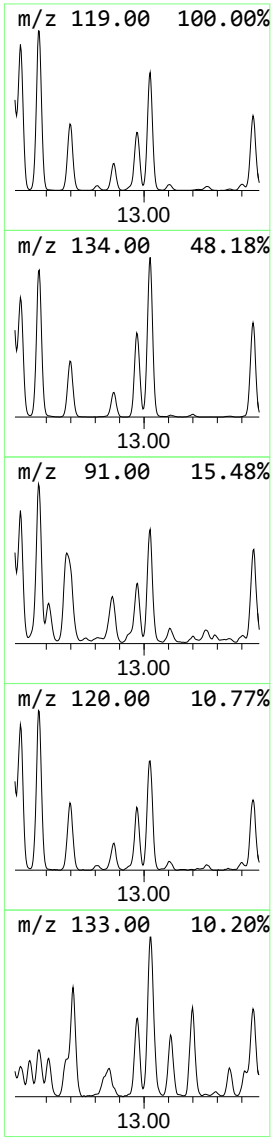
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.970	15.04 ug/L	533264	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
4	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

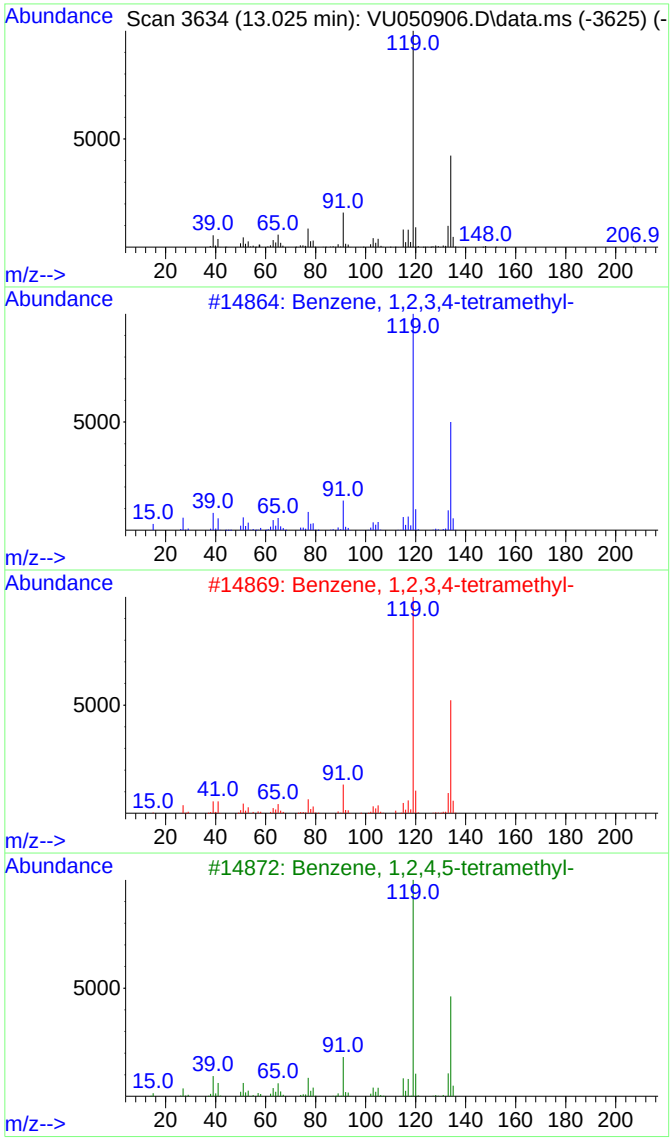
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

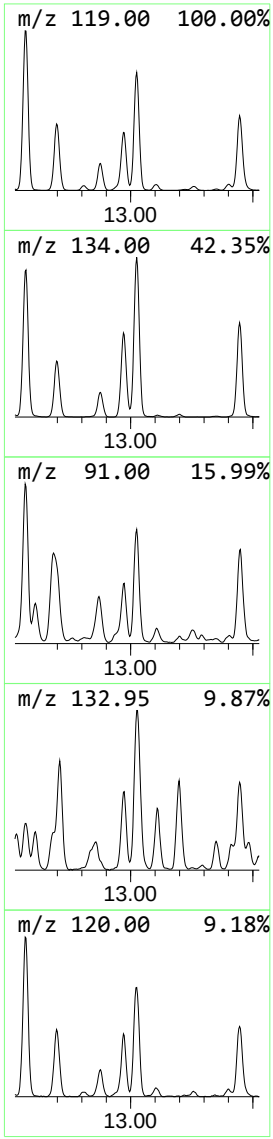
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.025	35.27 ug/L	1250390	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
4	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

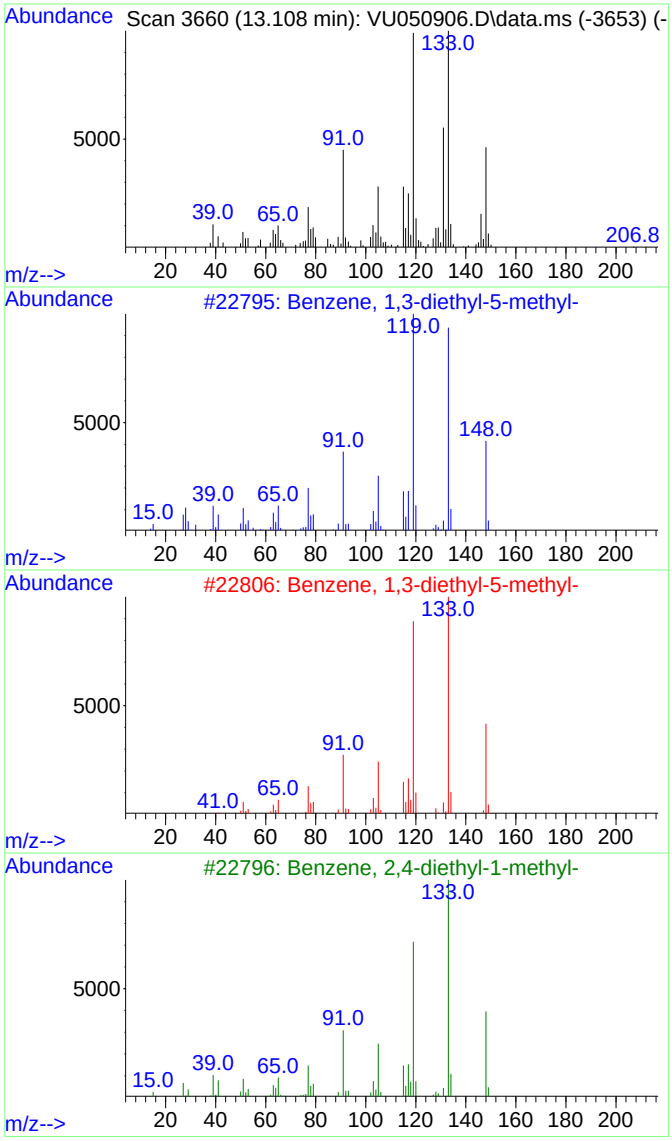
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

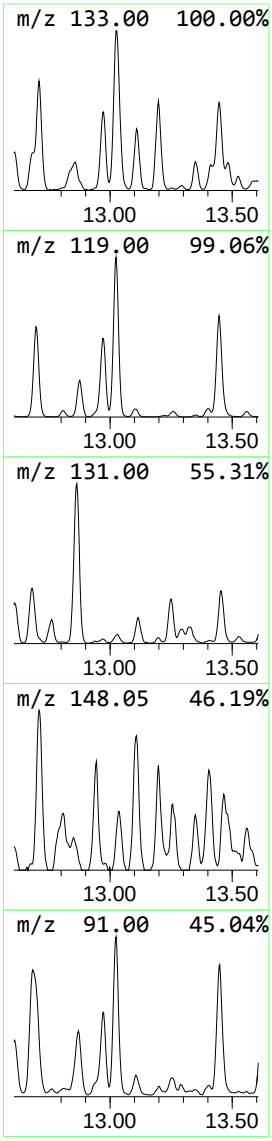
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 Benzene, 1,3-diethyl-5-methyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.108	4.13 ug/L	146523	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	96
2	Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	90
3	Benzene, 2,4-diethyl-1-methyl-	148	C11H16	001758-85-6	89
4	3,4-Dimethylcumene	148	C11H16	1000370-34-1	50
5	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	41





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

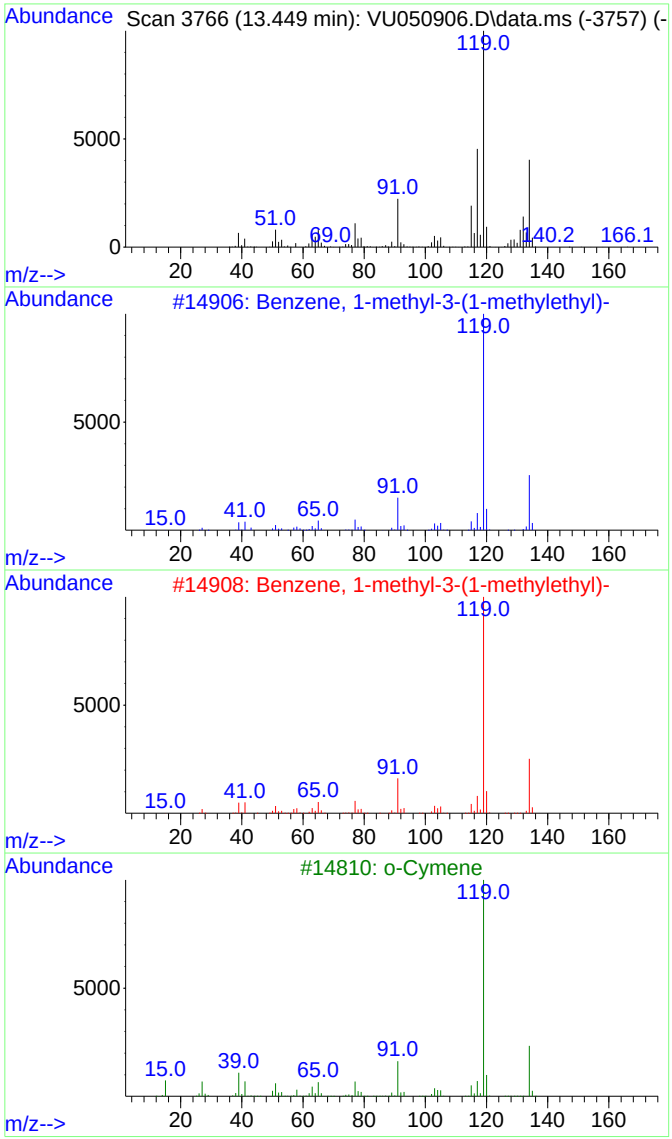
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

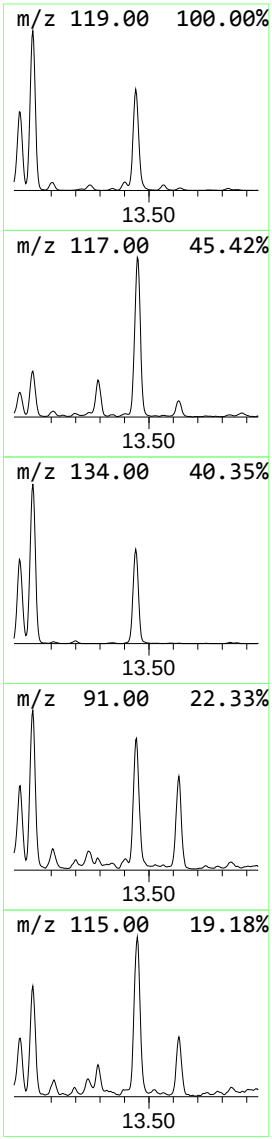
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 o-Cymene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.449	31.77 ug/L	1126190	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	64
2	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	64
3	o-Cymene	134	C10H14	000527-84-4	64
4	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	64
5	o-Cymene	134	C10H14	000527-84-4	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

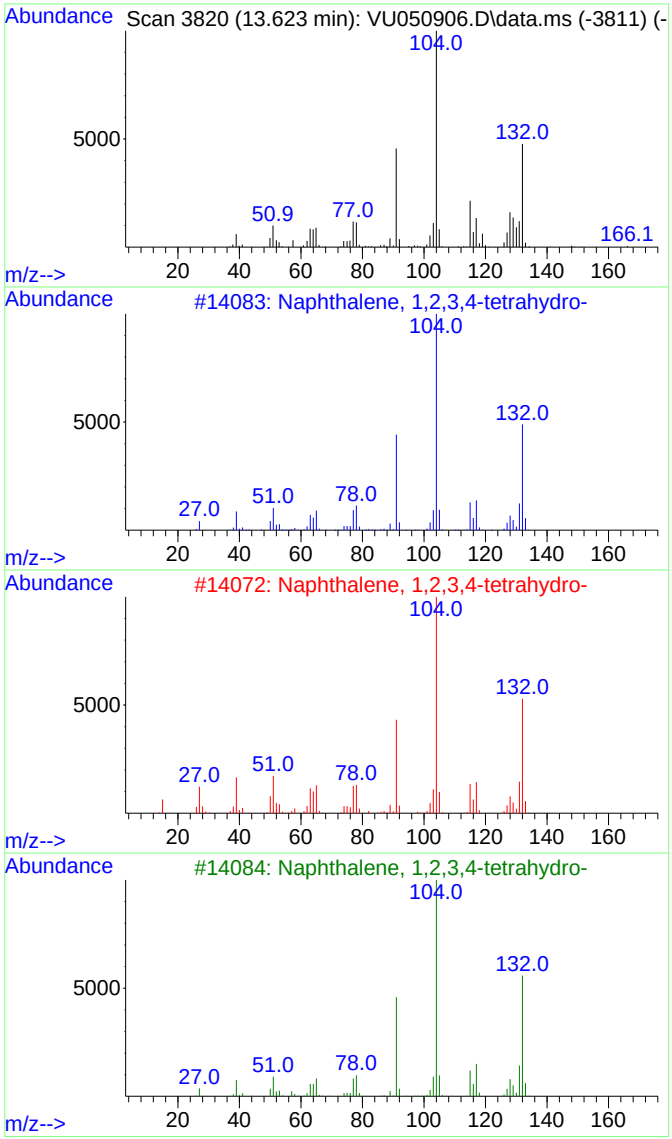
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

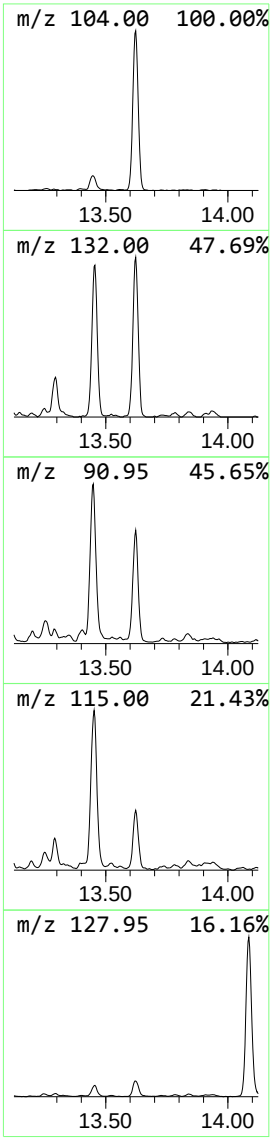
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.623	11.66 ug/L	413546	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	90
2	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	90
3	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	90
4	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	89
5	Indan, epoxide	132	C9H8O	000768-22-9	46





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

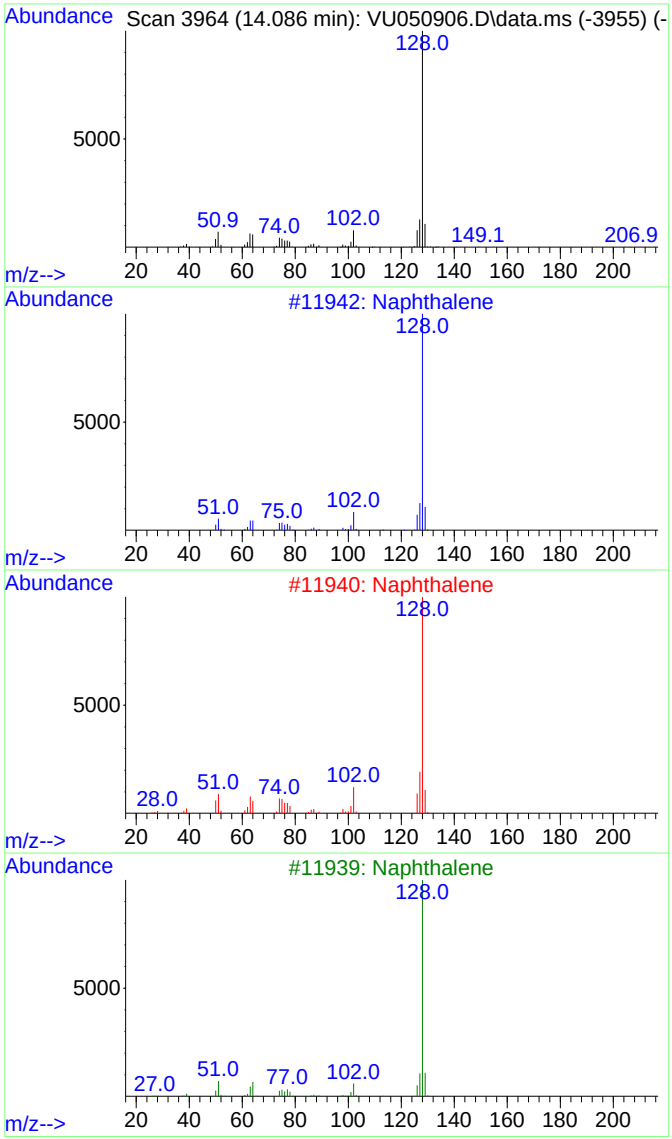
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

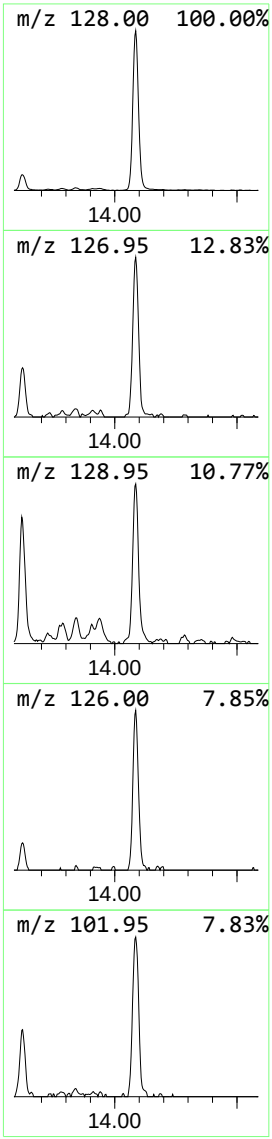
****

Peak Number 26 Azulene Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.086	9.36 ug/L	331673	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene	128	C10H8	000091-20-3	95
2	Naphthalene	128	C10H8	000091-20-3	94
3	Naphthalene	128	C10H8	000091-20-3	93
4	Azulene	128	C10H8	000275-51-4	91
5	1H-Indene, 1-methylene-	128	C10H8	002471-84-3	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092222\
Data File : VU050906.D
Acq On : 22 Sep 2022 18:25
Operator : JC/MD
Sample : N4747-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E45A3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092022WMA.M
Quant Title : VOC Analysis

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: C...	8.864	2.9	ug/L	69161	2	9.417	1204310	50.0
(DEL) Alkane: C...	8.922	3.3	ug/L	79820	2	9.417	1204310	50.0
Cyclohexene, 1-...	9.883	3.1	ug/L	74409	2	9.417	1204310	50.0
(DEL) Alkane: C...	10.034	8.7	ug/L	208405	2	9.417	1204310	50.0
Methyl 10,11-te...	10.269	16.1	ug/L	387279	2	9.417	1204310	50.0
(DEL) Alkane: C...	10.356	5.4	ug/L	129207	2	9.417	1204310	50.0
(DEL) Alkane: C...	10.809	8.8	ug/L	313165	3	11.809	1772660	50.0
Benzene, propyl-	10.906	27.9	ug/L	988524	3	11.809	1772660	50.0
Benzene, 1-ethy...	11.022	53.6	ug/L	1901050	3	11.809	1772660	50.0
Benzene, 1-ethy...	11.292	76.6	ug/L	2715270	3	11.809	1772660	50.0
Benzene, 1,2,3-...	11.893	125.7	ug/L	4456910	3	11.809	1772660	50.0
p-Cymene	12.005	9.6	ug/L	338579	3	11.809	1772660	50.0
Benzene, 1,4-di...	12.095	44.3	ug/L	1569320	3	11.809	1772660	50.0
Benzene, 1,2-di...	12.301	11.0	ug/L	391259	3	11.809	1772660	50.0
Benzene, 1-meth...	12.375	14.0	ug/L	495287	3	11.809	1772660	50.0
Benzene, 2-ethy...	12.462	32.7	ug/L	1159810	3	11.809	1772660	50.0
Benzene, 1-meth...	12.568	37.0	ug/L	1311100	3	11.809	1772660	50.0
Benzene, (2-met...	12.684	43.0	ug/L	1523950	3	11.809	1772660	50.0
Benzene, 4-ethy...	12.870	11.7	ug/L	415587	3	11.809	1772660	50.0
Benzene, 1,2,4,...	12.970	15.0	ug/L	533264	3	11.809	1772660	50.0
Benzene, 1,2,3,...	13.025	35.3	ug/L	1250390	3	11.809	1772660	50.0
Benzene, 1,3-di...	13.108	4.1	ug/L	146523	3	11.809	1772660	50.0
o-Cymene	13.449	31.8	ug/L	1126190	3	11.809	1772660	50.0
Naphthalene, 1,...	13.623	11.7	ug/L	413546	3	11.809	1772660	50.0
Azulene	14.086	9.4	ug/L	331673	3	11.809	1772660	50.0