LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028102.D
Acq On : 14 Apr 2022 13:17
Operator : JC/MD
Sample : N2403-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-24

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\82X041222W.M
Title : SW846 8260

Signal : TIC: VX028102.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.264	26	29	37	rBV	166888	149379	2.38%	0.249%
2	1.368	43	46	51	rBV	876254	831053	13.22%	1.384%
3	1.746	103	108	123	rVB	4765243	6284871	100.00%	10.469%
4	1.953	137	142	155	rVB2	855256	1301889	20.71%	2.169%
5	2.069	156	161	166	rBV	69250	94514	1.50%	0.157%
6	2.215	179	185	198	rVB	511103	890228	14.16%	1.483%
7	2.337	198	205	217	rVB	155826	325277	5.18%	0.542%
8	2.782	259	278	281	rBV2	435931	1064303	16.93%	1.773%
9	2.825	281	285	288	rVV	592460	1165591	18.55%	1.942%
10	2.867	288	292	303	rVV2	714061	1581372	25.16%	2.634%
11	2.953	303	306	320	rVB	40245	87054	1.39%	0.145%
12	3.105	320	331	346	rBV2	635560	1469026	23.37%	2.447%
13	3.318	357	366	377	rBV	56252	126832	2.02%	0.211%
14	3.575	400	408	418	rVB	28894	69073	1.10%	0.115%
15	3.733	423	434	443	rBV	282327	707607	11.26%	1.179%
16	3.837	443	451	459	rBV	199581	494924	7.87%	0.824%
17	3.995	469	477	482	rBV3	35412	93219	1.48%	0.155%
18	4.105	484	495	504	rVV	190336	535574	8.52%	0.892%
19	4.215	504	513	518	rVV2	79902	232535	3.70%	0.387%
20	4.306	518	528	539	rVV	1048373	2960715	47.11%	4.932%
21	4.428	539	548	562	rVB	141282	426385	6.78%	0.710%
22	5.196	649	674	689	rVB	310480	1069592	17.02%	1.782%
23	5.385	693	705	711	rBV2	173839	516726	8.22%	0.861%
24	5.477	711	720	728	rVV	658558	2185463	34.77%	3.641%
25	5.556	728	733	739	rVV	259227	759468	12.08%	1.265%
26	5.623	739	744	765	rVB2	220412	725215	11.54%	1.208%
27	5.830	765	778	791	rBV4	144540	474777	7.55%	0.791%
28	5.958	791	799	804	rVV2	163953	425487	6.77%	0.709%
29	6.044	804	813	824	rVV	1800395	4716366	75.04%	7.857%
30	6.166	824	833	835	rVV2	146352	372491	5.93%	0.620%
31	6.251	836	847	857	rVV5	407177	1740094	27.69%	2.899%
32	6.361	857	865	877	rVB3	165772	473058	7.53%	0.788%
33	6.617	892	907	921	rVB4	32300	129630	2.06%	0.216%
34	6.763	921	931	941	rBV2	425250	1293072	20.57%	2.154%
35	6.854	942	946	958	rVB2	186752	465278	7.40%	0.775%
36	6.976	958	966	973	rBV3	38551	90959	1.45%	0.152%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028102.D
Acq On : 14 Apr 2022 13:17
Operator : JC/MD
Sample : N2403-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-24

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\82X041222W.M
Title : SW846 8260

37	7.068	973	981	989	rVB4	29438	67305	1.07%	0.112%
38	7.171	989	998	1007	rBV2	206846	460104	7.32%	0.766%
39	7.379	1019	1032	1041	rBV2	565024	1468527	23.37%	2.446%
40	7.659	1073	1078	1090	rVB	65178	126845	2.02%	0.211%
41	7.775	1090	1097	1104	rVB2	38404	78893	1.26%	0.131%
42	7.860	1104	1111	1112	rBV2	78648	136468	2.17%	0.227%
43	7.885	1112	1115	1122	rVV3	101742	209966	3.34%	0.350%
44	7.988	1122	1132	1143	rVB2	150350	402477	6.40%	0.670%
45	8.110	1143	1152	1154	rBV	246708	468062	7.45%	0.780%
46	8.147	1154	1158	1162	rVV	301244	542262	8.63%	0.903%
47	8.196	1162	1166	1176	rVB4	74630	156760	2.49%	0.261%
48	8.311	1176	1185	1192	rBV5	48701	115922	1.84%	0.193%
49	8.427	1199	1204	1209	rBV	86285	139987	2.23%	0.233%
50	8.507	1209	1217	1224	rVB	362776	623205	9.92%	1.038%
51	8.610	1224	1234	1235	rBV4	93573	193574	3.08%	0.322%
52	8.653	1235	1241	1247	rVV	856811	1441404	22.93%	2.401%
53	8.720	1247	1252	1257	rVB2	84401	136927	2.18%	0.228%
54	8.842	1268	1272	1279	rVB	104704	184876	2.94%	0.308%
55	8.927	1281	1286	1289	rBV	53243	88500	1.41%	0.147%
56	8.976	1289	1294	1299	rVB3	57448	109968	1.75%	0.183%
57	9.104	1311	1315	1320	rVB	52844	84073	1.34%	0.140%
58	9.165	1320	1325	1334	rVB	180034	302138	4.81%	0.503%
59	9.378	1356	1360	1364	rBV	68548	94844	1.51%	0.158%
60	9.458	1364	1373	1377	rBV	472943	757024	12.05%	1.261%
61	9.500	1377	1380	1386	rVB4	54396	96681	1.54%	0.161%
62	9.561	1386	1390	1394	rBV	70337	105300	1.68%	0.175%
63	10.055	1460	1471	1476	rBV	961483	1334396	21.23%	2.223%
64	10.147	1482	1486	1490	rVB3	42011	71440	1.14%	0.119%
65	10.195	1490	1494	1499	rVV2	175936	320154	5.09%	0.533%
66	10.244	1499	1502	1508	rVV	144518	244631	3.89%	0.408%
67	10.305	1508	1512	1518	rVB	403508	522002	8.31%	0.870%
68	10.415	1526	1530	1534	rVB2	118276	162060	2.58%	0.270%
69	10.646	1564	1568	1573	rVB2	61334	85607	1.36%	0.143%
70	10.872	1598	1605	1614	rBV3	162154	338366	5.38%	0.564%
71	10.963	1614	1620	1625	rVB	240658	335826	5.34%	0.559%
72	11.085	1635	1640	1645	rVV	751655	985586	15.68%	1.642%
73	11.171	1649	1654	1660	rVB	241861	330678	5.26%	0.551%
74	11.305	1672	1676	1684	rVB	313828	385084	6.13%	0.641%
75	11.402	1685	1692	1695	rBV	99099	133228	2.12%	0.222%
76	11.494	1703	1707	1712	rVB	139684	172899	2.75%	0.288%

LSC Area Percent Report

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Data File : VX028102.D
Acq On : 14 Apr 2022 13:17
Operator : JC/MD
Sample : N2403-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-24

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\82X041222W.M
Title : SW846 8260

77	11.616	1721	1727	1734	rBV	498715	641930	10.21%	1.069%
78	11.707	1738	1742	1746	rVV	191041	270675	4.31%	0.451%
79	11.750	1746	1749	1755	rVB	99761	131216	2.09%	0.219%
80	12.024	1789	1794	1800	rBV	1021656	1246329	19.83%	2.076%
81	12.085	1800	1804	1809	rBV	136616	184380	2.93%	0.307%
82	12.244	1825	1830	1835	rBV	2388850	2867740	45.63%	4.777%
83	12.317	1839	1842	1849	rVB2	71618	114994	1.83%	0.192%
84	12.408	1853	1857	1861	rVV2	65312	84398	1.34%	0.141%
85	12.463	1861	1866	1872	rVB	111596	149878	2.38%	0.250%
86	12.530	1872	1877	1885	rBV2	78497	146719	2.33%	0.244%
87	12.609	1885	1890	1894	rVV2	330402	476453	7.58%	0.794%
88	12.646	1894	1896	1900	rVV	113606	136642	2.17%	0.228%
89	12.701	1900	1905	1918	rVB	479721	712050	11.33%	1.186%
90	12.847	1925	1929	1934	rVB2	50826	70464	1.12%	0.117%
91	12.920	1934	1941	1945	rBV2	155721	229789	3.66%	0.383%
92	12.963	1945	1948	1954	rVB	74716	97010	1.54%	0.162%
93	13.036	1954	1960	1965	rBV4	43009	92608	1.47%	0.154%
94	13.170	1979	1982	1985	rBV	162322	169987	2.70%	0.283%
95	13.292	1998	2002	2006	rVV2	467867	610442	9.71%	1.017%
96	13.341	2006	2010	2020	rVV5	117542	199938	3.18%	0.333%
97	13.426	2020	2024	2029	rVB2	119170	155453	2.47%	0.259%
98	13.664	2061	2063	2069	rVB2	56980	76991	1.23%	0.128%
99	13.780	2076	2082	2090	rVB	540110	706139	11.24%	1.176%
100	14.780	2241	2246	2253	rBV2	79744	115970	1.85%	0.193%

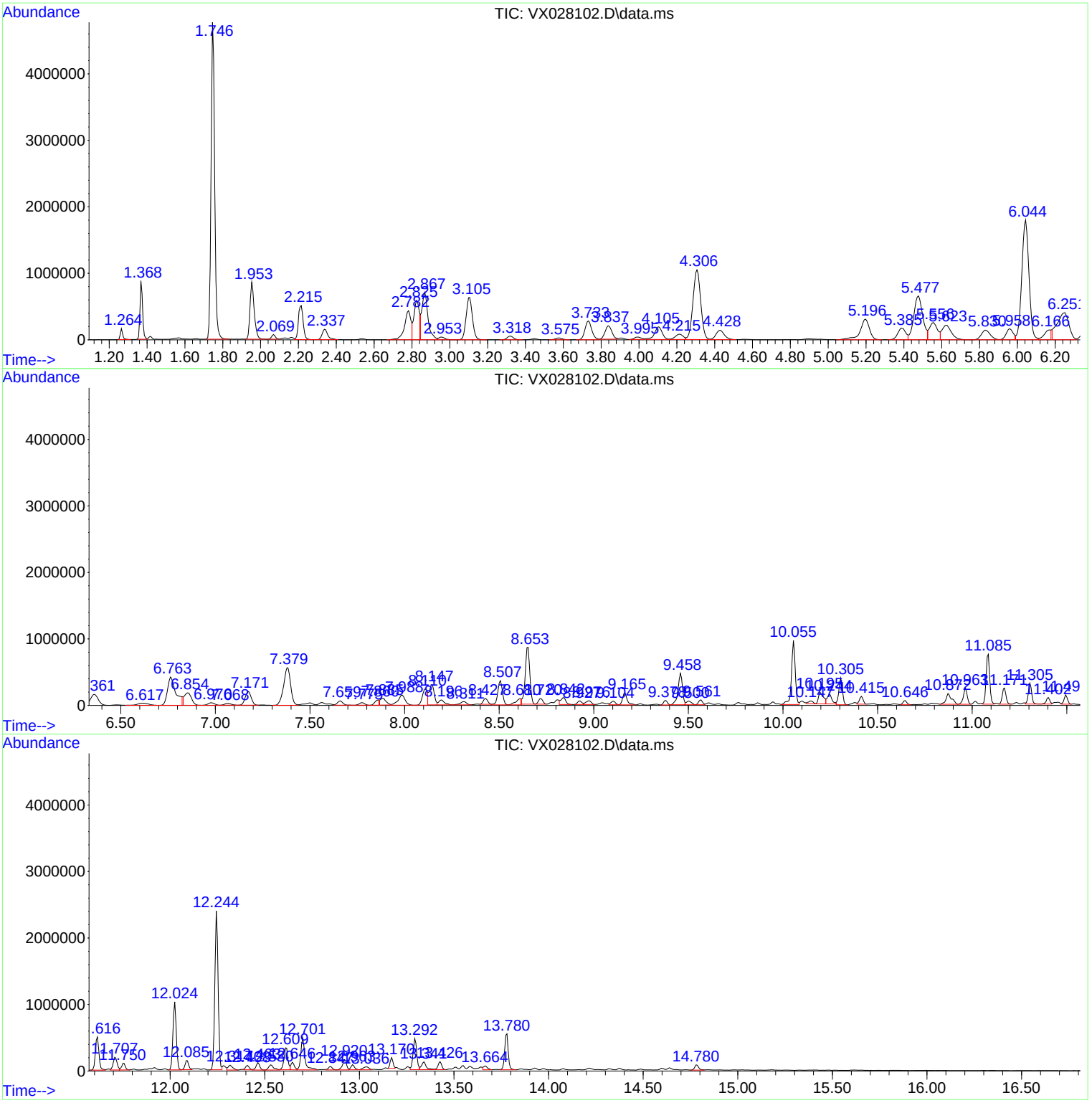
Sum of corrected areas: 60031341

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028102.D
Acq On : 14 Apr 2022 13:17
Operator : JC/MD
Sample : N2403-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-24

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028102.D
Acq On : 14 Apr 2022 13:17
Operator : JC/MD
Sample : N2403-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-24

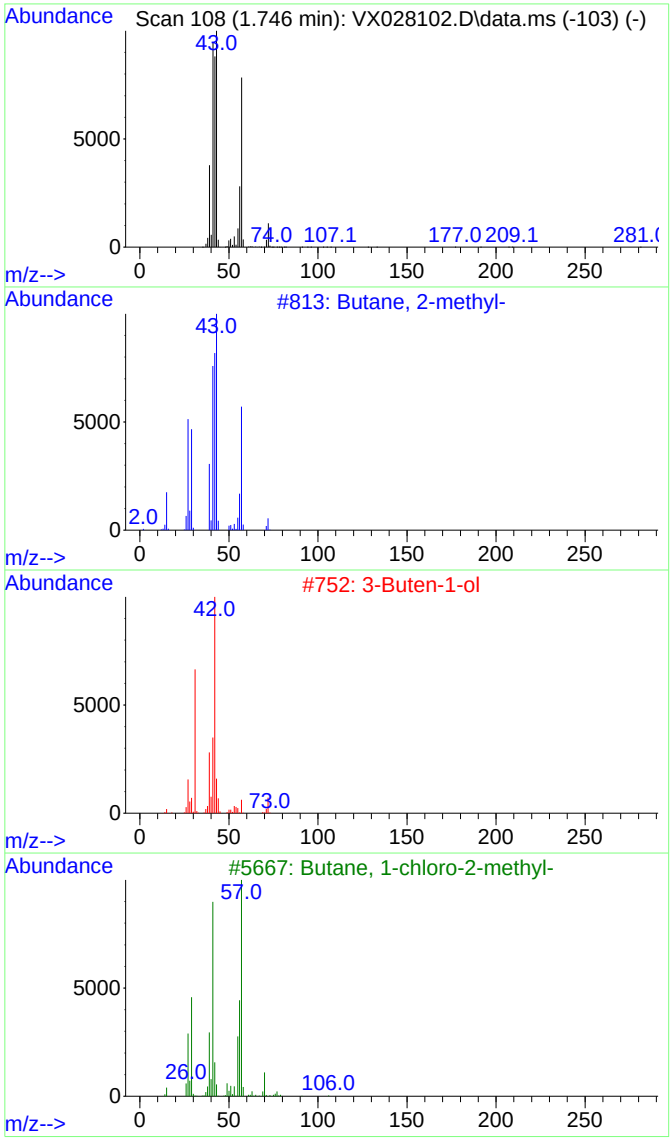
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

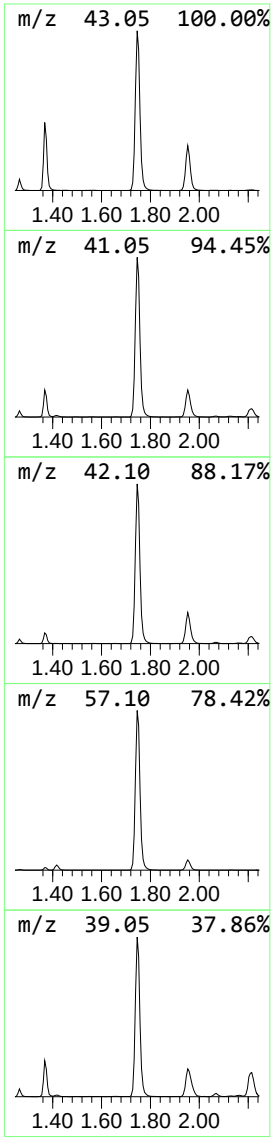
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Butane, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.746	413.77 ug/l	6284870	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	86
2	3-Buten-1-ol	72	C4H8O	000627-27-0	43
3	Butane, 1-chloro-2-methyl-	106	C5H11Cl	000616-13-7	35
4	Oxirane, trimethyl-	86	C5H10O	005076-19-7	33
5	Pentane	72	C5H12	000109-66-0	28





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028102.D
Acq On : 14 Apr 2022 13:17
Operator : JC/MD
Sample : N2403-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-24

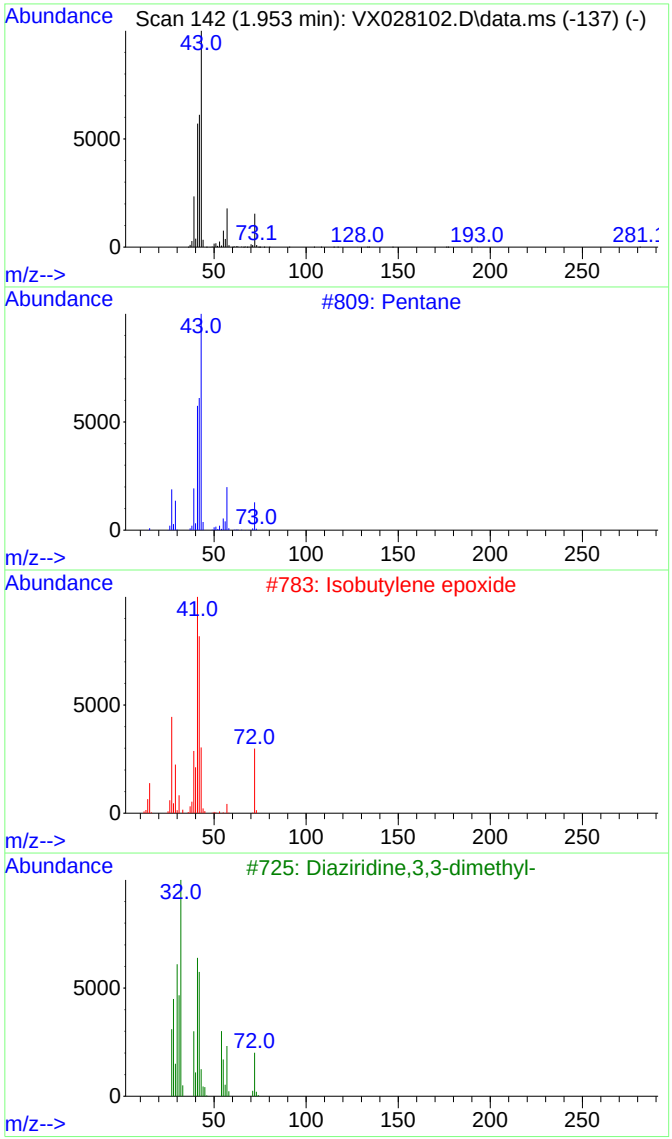
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

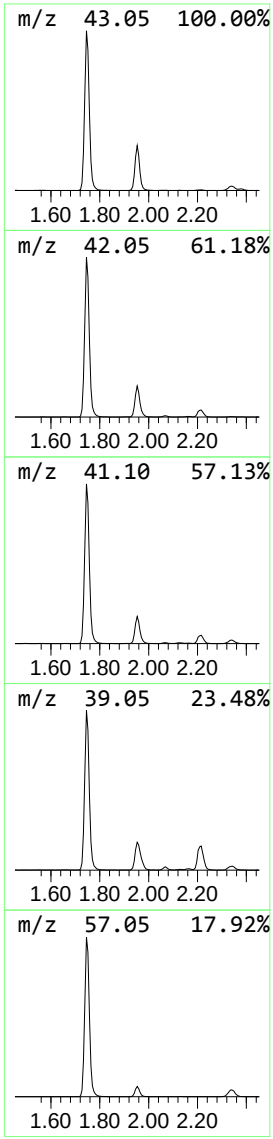
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Pentane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.953	85.71 ug/l	1301890	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane	72	C5H12	000109-66-0	91
2	Isobutylene epoxide	72	C4H8O	000558-30-5	47
3	Diaziridine,3,3-dimethyl-	72	C3H8N2	004901-76-2	42
4	Oxirane, ethyl-	72	C4H8O	000106-88-7	40
5	1-Propene, 2-methoxy-	72	C4H8O	000116-11-0	37





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028102.D
Acq On : 14 Apr 2022 13:17
Operator : JC/MD
Sample : N2403-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-24

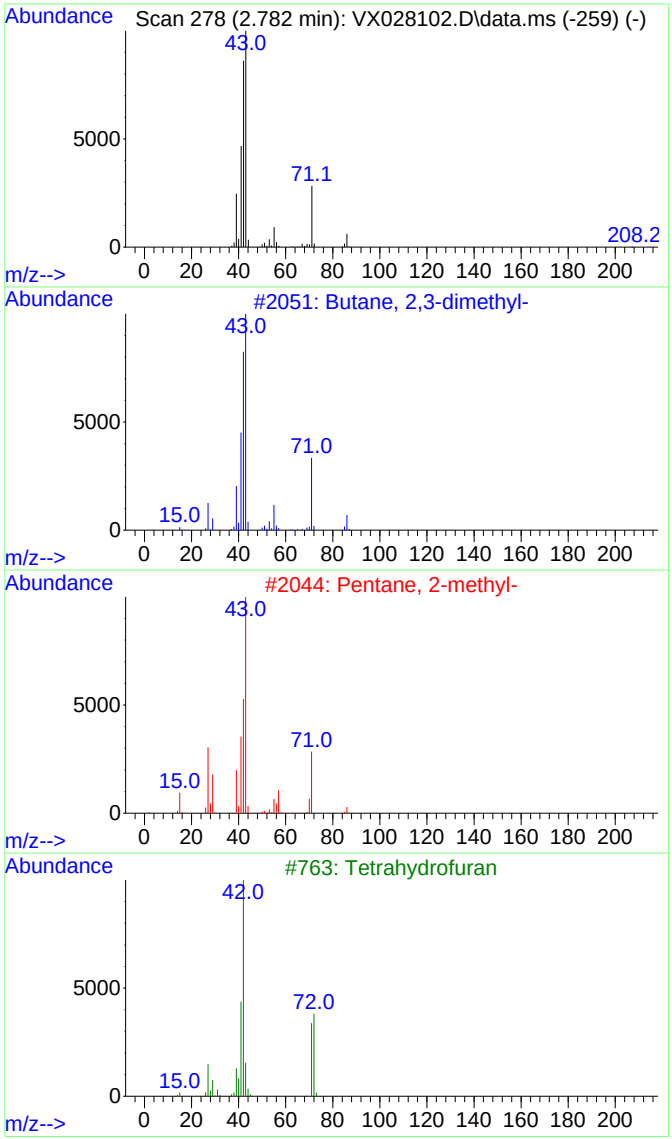
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

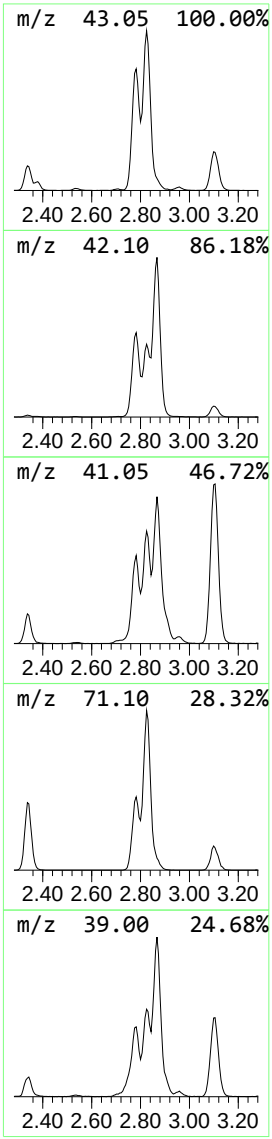
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Butane, 2,3-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.782	70.07 ug/l	1064300	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	91
2	Pentane, 2-methyl-	86	C6H14	000107-83-5	50
3	Tetrahydrofuran	72	C4H8O	000109-99-9	38
4	1-Pentene	70	C5H10	000109-67-1	28
5	Pyrrolidine	71	C4H9N	000123-75-1	12





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028102.D
Acq On : 14 Apr 2022 13:17
Operator : JC/MD
Sample : N2403-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-24

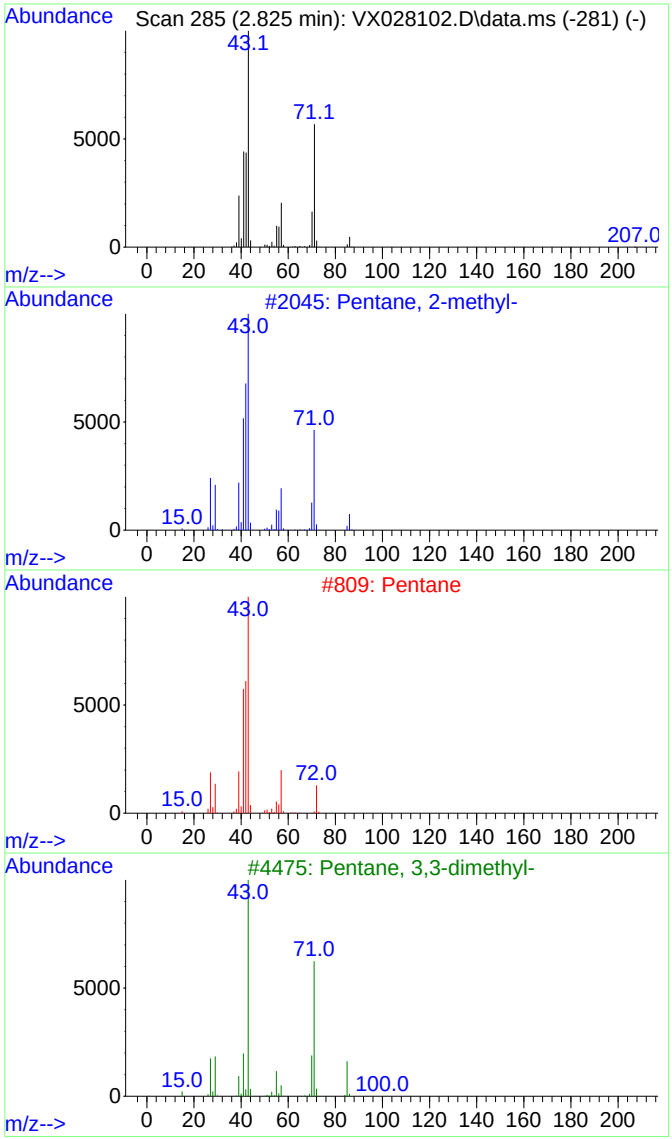
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

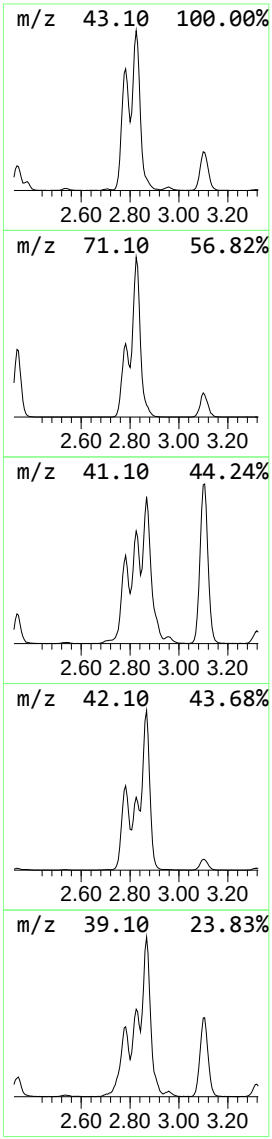
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Pentane, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.825	76.74 ug/l	1165590	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2-methyl-	86	C6H14	000107-83-5	94
2	Pentane	72	C5H12	000109-66-0	43
3	Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	43
4	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	38
5	1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028102.D
Acq On : 14 Apr 2022 13:17
Operator : JC/MD
Sample : N2403-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-24

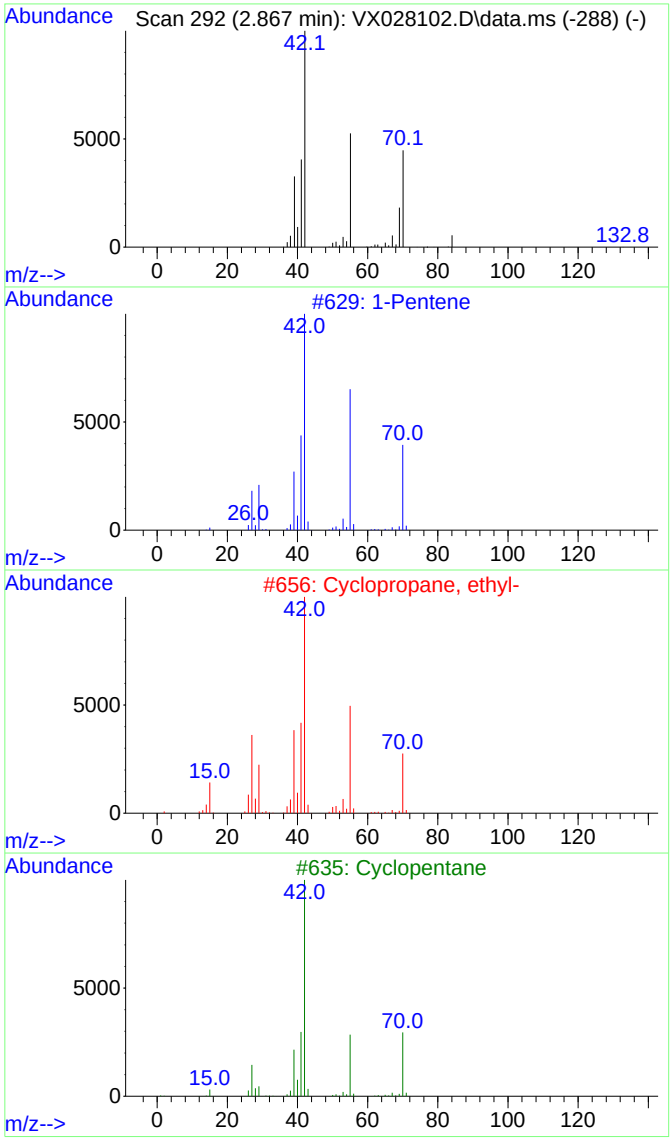
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

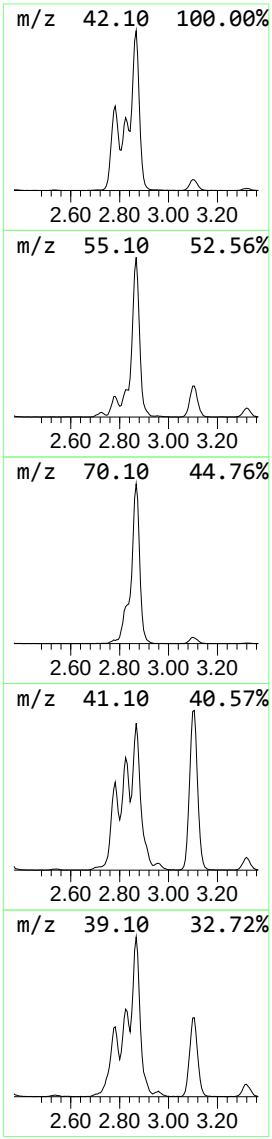
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 1-Pentene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.867	104.11 ug/l	1581370	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentene	70	C5H10	000109-67-1	80
2	Cyclopropane, ethyl-	70	C5H10	001191-96-4	59
3	Cyclopentane	70	C5H10	000287-92-3	56
4	2-Pentene, (E)-	70	C5H10	000646-04-8	43
5	3-Buten-1-ol	72	C4H8O	000627-27-0	37





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028102.D
Acq On : 14 Apr 2022 13:17
Operator : JC/MD
Sample : N2403-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-24

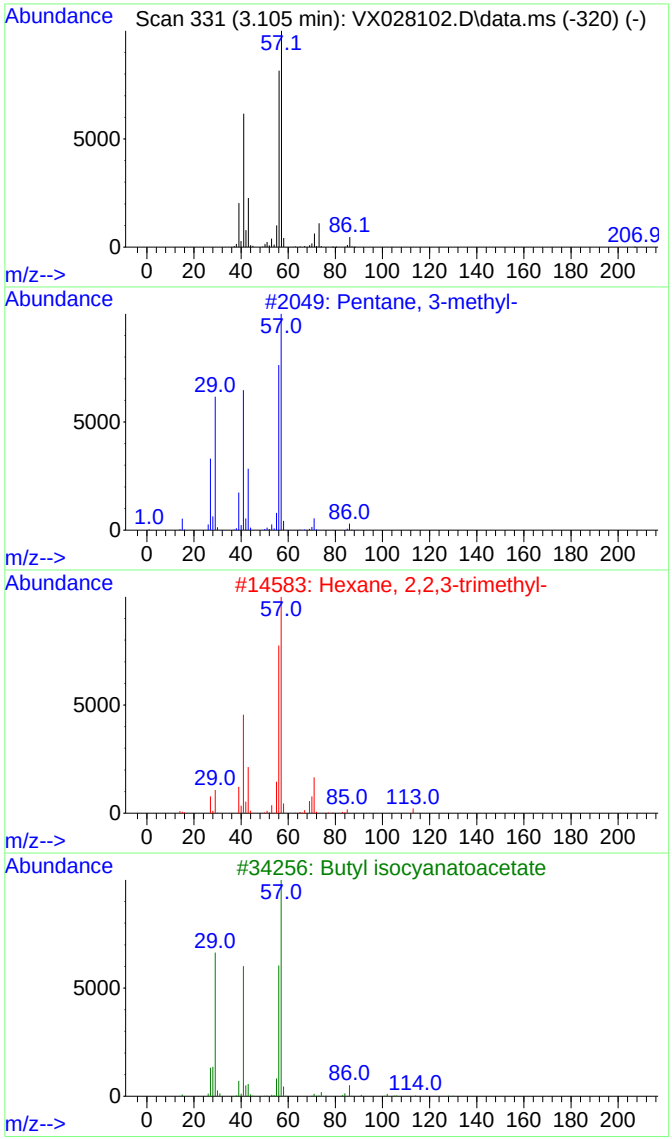
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

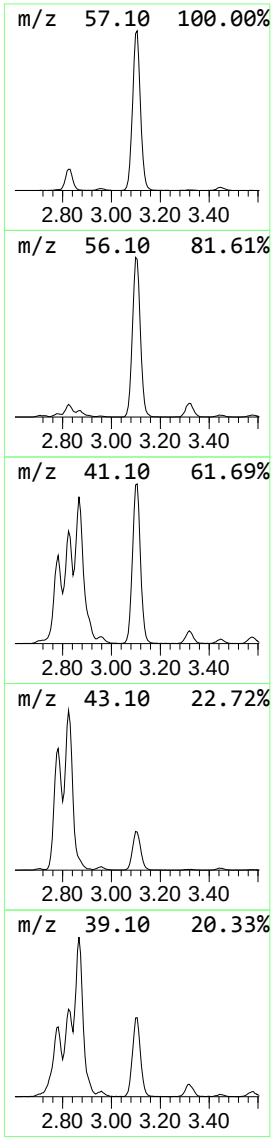
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Pentane, 3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.105	96.71 ug/l	1469030	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	87
2	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	72
3	Butyl isocyanatoacetate	157	C7H11NO3	017046-22-9	56
4	3,4-Dimethyldihydrofuran-2,5-dione	128	C6H8O3	007475-92-5	50
5	Sulphuric acid dibutyl ester	210	C8H18O4S	000625-22-9	39





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Data File : VX028102.D
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Operator : JC/MD
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-24

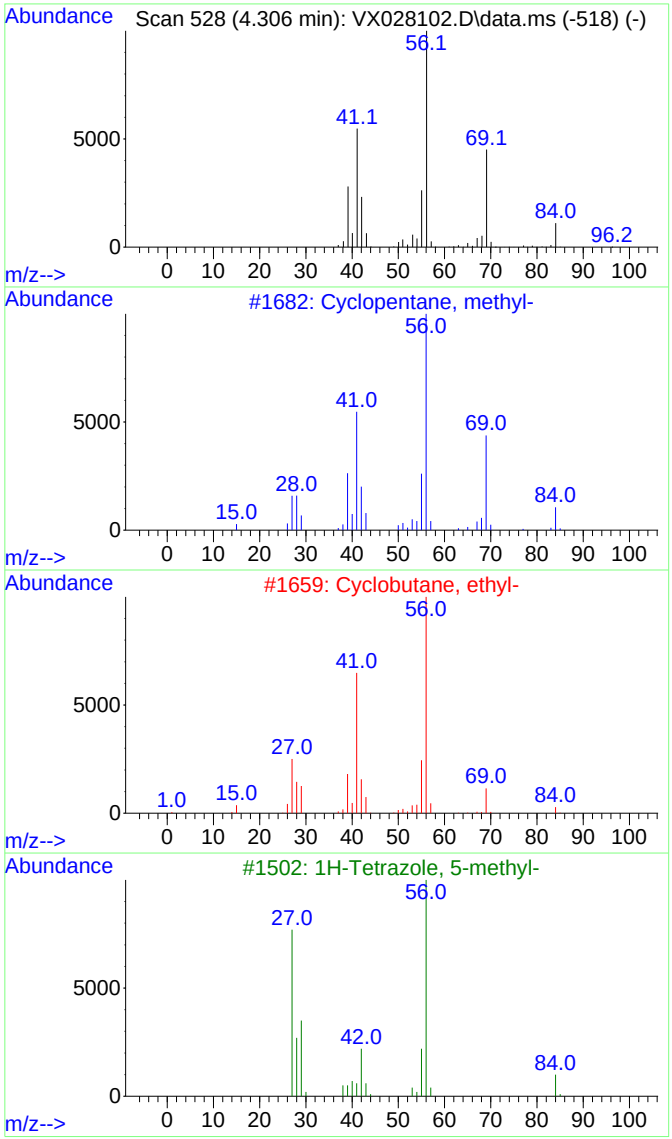
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

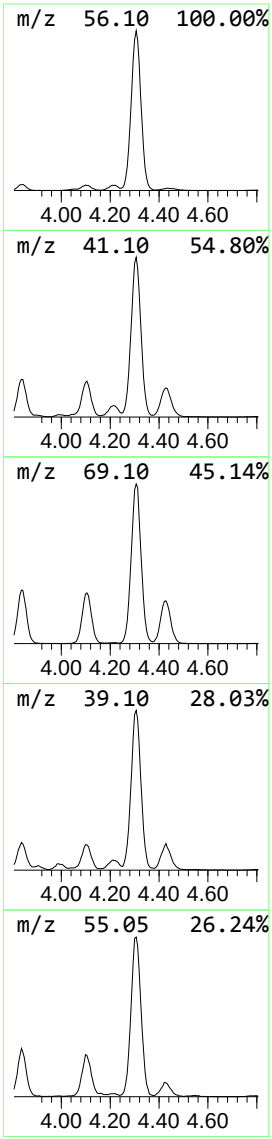
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Cyclopentane, methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.306	194.92 ug/l	2960720	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	87
2	Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
3	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	50
4	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	49
5	Cyclopropane, 1-ethyl-2-methyl-,...	84	C6H12	019781-68-1	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028102.D
Acq On : 14 Apr 2022 13:17
Operator : JC/MD
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-24

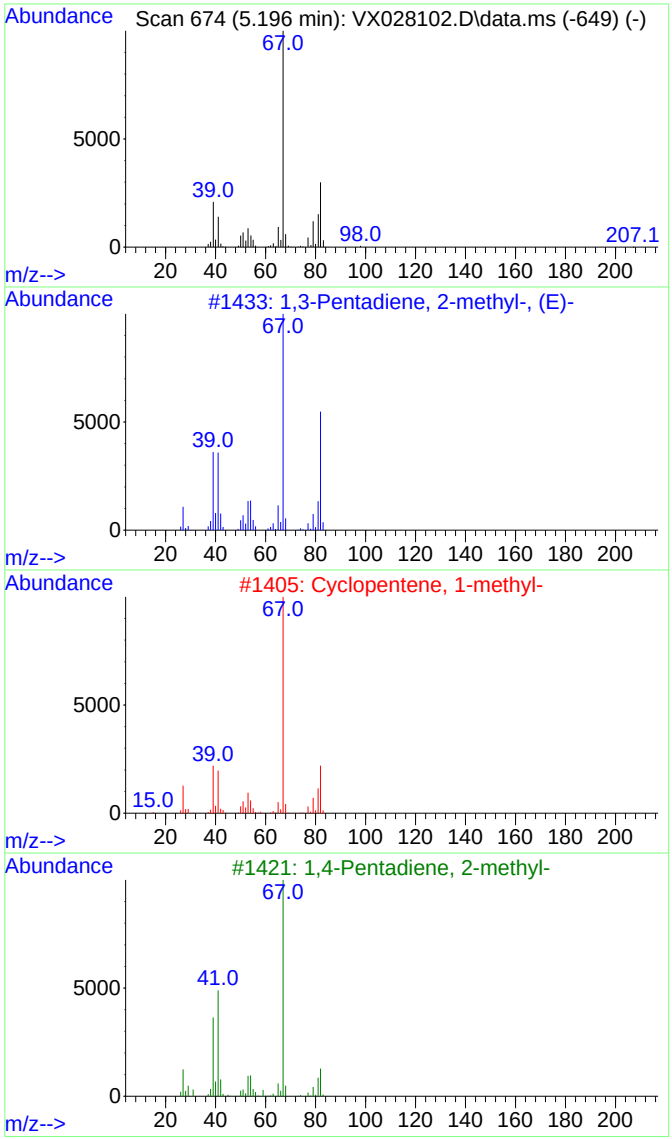
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

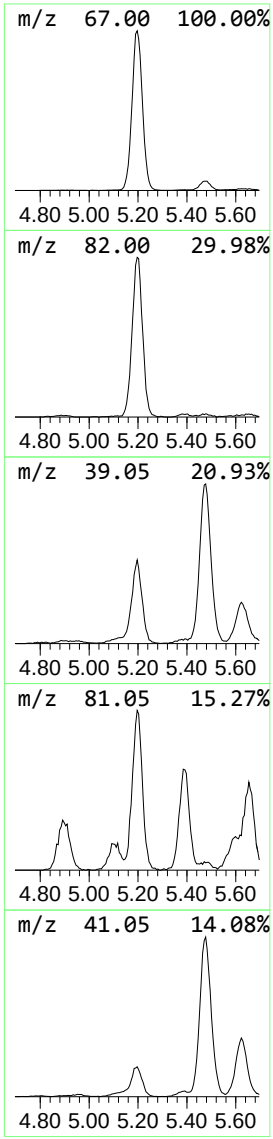
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 1,3-Pentadiene, 2-methyl-, ... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.196	70.42 ug/l	1069590	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Pentadiene, 2-methyl-, (E)-	82	C6H10	000926-54-5	90
2	Cyclopentene, 1-methyl-	82	C6H10	000693-89-0	90
3	1,4-Pentadiene, 2-methyl-	82	C6H10	000763-30-4	90
4	Cyclohexene	82	C6H10	000110-83-8	81
5	Cyclopentane, methylene-	82	C6H10	001528-30-9	81





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Operator : JC/MD
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ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-24

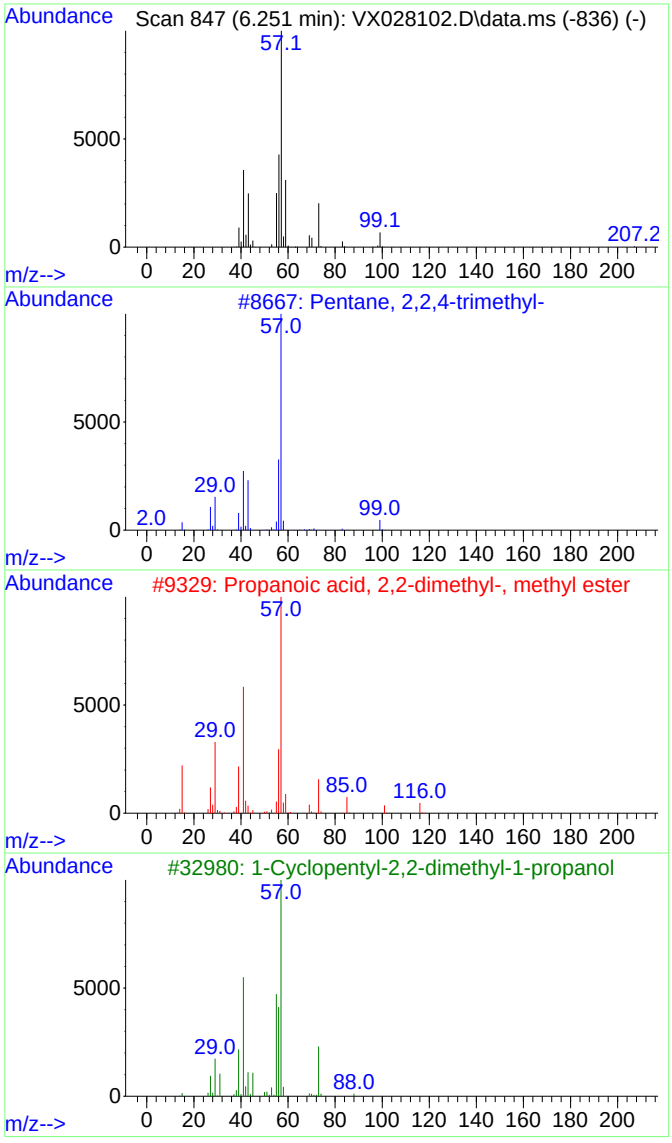
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

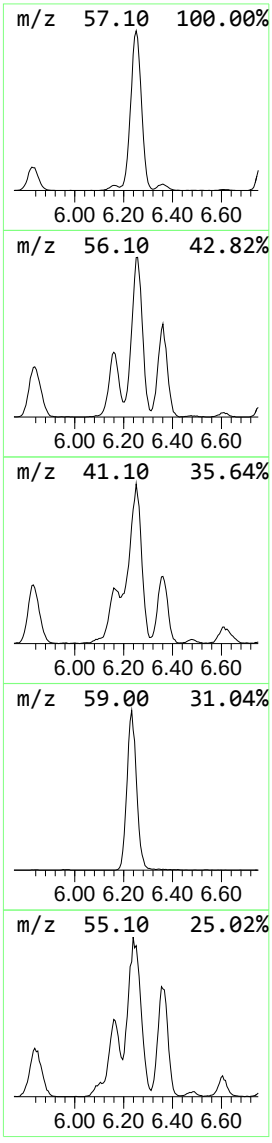
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 unknown6.251 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.251	67.29 ug/l	1740090	1,4-Difluorobenzene	6.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	47
2	Propanoic acid, 2,2-dimethyl-, m...	116	C6H12O2	000598-98-1	40
3	1-Cyclopentyl-2,2-dimethyl-1-pro...	156	C10H20O	337966-85-5	36
4	1-Propanol, 2,2-dimethyl-	88	C5H12O	000075-84-3	36
5	Dodecane, 2,2,11,11-tetramethyl-	226	C16H34	127204-12-0	33





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ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-24

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

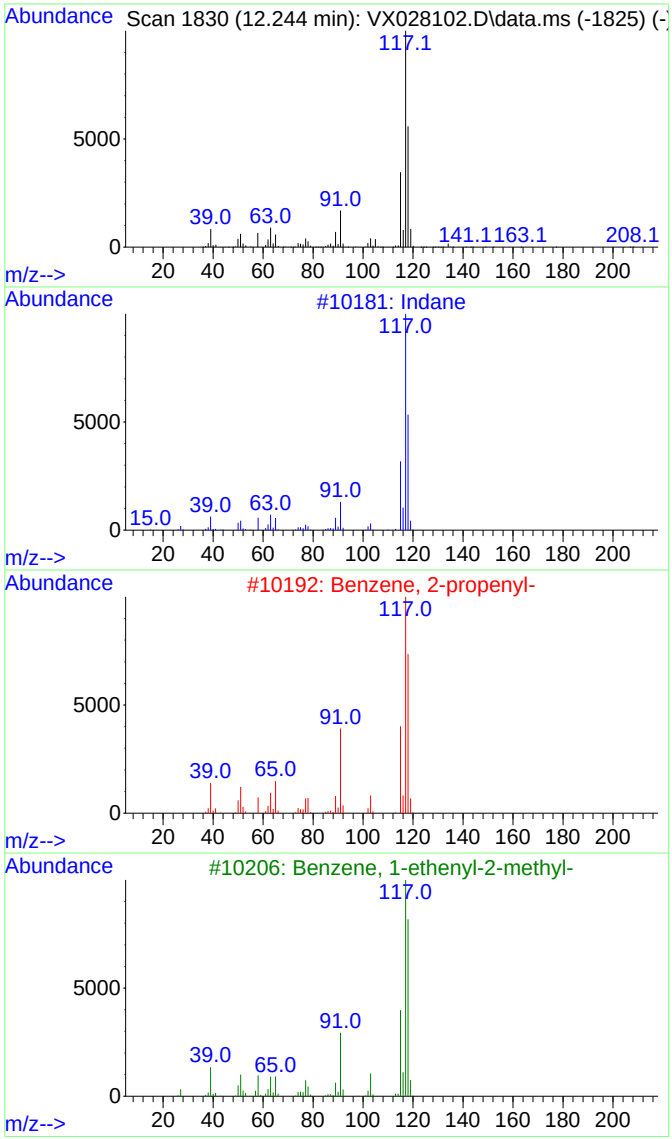
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

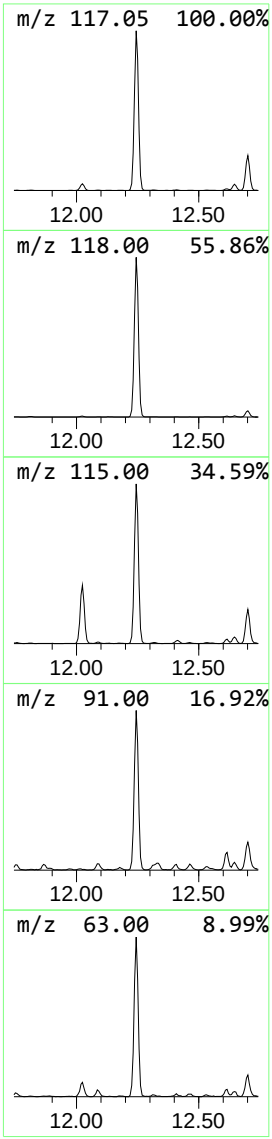
****

Peak Number 10 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	115.05 ug/l	2867740	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	91
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	74
3	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	74
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	72
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	64





Tentatively Identified Compound (LSC) summary

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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-methyl-	1.746	413.8	ug/l	6284870	1	5.556	759468	50.0
Pentane	1.953	85.7	ug/l	1301890	1	5.556	759468	50.0
Butane, 2,3-dim...	2.782	70.1	ug/l	1064300	1	5.556	759468	50.0
Pentane, 2-methyl-	2.825	76.7	ug/l	1165590	1	5.556	759468	50.0
1-Pentene	2.867	104.1	ug/l	1581370	1	5.556	759468	50.0
Pentane, 3-methyl-	3.105	96.7	ug/l	1469030	1	5.556	759468	50.0
Cyclopentane, m...	4.306	194.9	ug/l	2960720	1	5.556	759468	50.0
1,3-Pentadiene,...	5.196	70.4	ug/l	1069590	1	5.556	759468	50.0
unknown6.251	6.251	67.3	ug/l	1740090	2	6.763	1293070	50.0
Indane	12.244	115.1	ug/l	2867740	4	12.024	1246330	50.0