LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101123\
Data File : VX038225.D
Acq On : 11 Oct 2023 13:36
Operator : JC/MD
Sample : 04788-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

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\82X100523W.M
Title : SW846 8260

Signal : TIC: VX038225.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.374	43	47	52	rBV	69458	71330	2.07%	0.197%
2	1.752	104	109	121	rVB	552215	714962	20.72%	1.977%
3	1.959	138	143	154	rVB	356970	521198	15.10%	1.441%
4	2.069	154	161	167	rBV	47918	68817	1.99%	0.190%
5	2.215	181	185	195	rVB	62728	86982	2.52%	0.240%
6	2.343	198	206	210	rBV	29082	58493	1.70%	0.162%
7	2.715	258	267	272	rBV4	21990	59175	1.71%	0.164%
8	2.782	272	278	281	rBV	91723	185215	5.37%	0.512%
9	2.825	281	285	290	rVV	198394	352486	10.21%	0.974%
10	3.105	321	331	346	rBV	251818	569621	16.51%	1.575%
11	3.325	357	367	378	rBV2	52228	138147	4.00%	0.382%
12	3.446	378	387	400	rBV	148462	332603	9.64%	0.920%
13	3.587	400	410	415	rBV4	18626	47332	1.37%	0.131%
14	3.672	415	424	430	rVV	52894	135293	3.92%	0.374%
15	3.733	430	434	442	rVB3	15708	35839	1.04%	0.099%
16	3.837	442	451	457	rBV2	36600	93188	2.70%	0.258%
17	3.904	457	462	465	rVV	23036	53350	1.55%	0.147%
18	3.946	466	469	484	rVV3	29393	94607	2.74%	0.262%
19	4.105	484	495	503	rVB4	28588	76015	2.20%	0.210%
20	4.306	518	528	543	rVB	470658	1341379	38.87%	3.708%
21	5.196	667	674	687	rVB2	32502	107724	3.12%	0.298%
22	5.385	690	705	711	rBV	85965	266770	7.73%	0.738%
23	5.477	711	720	727	rVV	267675	786910	22.80%	2.175%
24	5.556	728	733	739	rVB	77891	186011	5.39%	0.514%
25	5.836	765	779	790	rBV5	67550	235917	6.84%	0.652%
26	5.958	790	799	814	rVB	84555	225305	6.53%	0.623%
27	6.159	814	832	842	rBV4	91288	333507	9.66%	0.922%
28	6.263	842	849	856	rVV	83210	235250	6.82%	0.650%
29	6.361	856	865	875	rVV	106763	312088	9.04%	0.863%
30	6.611	897	906	913	rBV4	22256	55358	1.60%	0.153%
31	6.763	923	931	938	rBV	199324	506241	14.67%	1.400%
32	7.068	974	981	990	rVB	19600	45402	1.32%	0.126%
33	7.177	990	999	1008	rBV4	36134	90565	2.62%	0.250%
34	7.379	1019	1032	1045	rBV	549488	1353374	39.22%	3.741%
35	7.659	1070	1078	1091	rVB	59997	120357	3.49%	0.333%
36	7.854	1103	1110	1111	rBV2	31973	46545	1.35%	0.129%

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MW-3

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

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Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.M
Title : SW846 8260

37	7.885	1112	1115	1121	rVV	36361	63234	1.83%	0.175%
38	8.318	1177	1186	1193	rBV6	16825	42448	1.23%	0.117%
39	8.500	1210	1216	1226	rVB2	82448	150025	4.35%	0.415%
40	8.598	1226	1232	1235	rBV2	42488	77176	2.24%	0.213%
41	8.647	1235	1240	1247	rVV	430123	717078	20.78%	1.982%
42	8.720	1247	1252	1257	rVV	134342	215566	6.25%	0.596%
43	8.775	1257	1261	1265	rVV	26095	47110	1.37%	0.130%
44	8.842	1265	1272	1281	rVB7	18586	52822	1.53%	0.146%
45	8.976	1289	1294	1300	rVB2	39847	62717	1.82%	0.173%
46	9.104	1305	1315	1320	rBV2	30428	59583	1.73%	0.165%
47	9.159	1320	1324	1334	rVB	53133	93805	2.72%	0.259%
48	9.433	1363	1369	1371	rBV2	24630	42689	1.24%	0.118%
49	9.561	1388	1390	1395	rVB2	31607	41187	1.19%	0.114%
50	10.055	1461	1471	1476	rBV	465196	646557	18.74%	1.787%
51	10.128	1476	1483	1488	rVB3	27944	55901	1.62%	0.155%
52	10.195	1488	1494	1499	rBV	190039	258155	7.48%	0.714%
53	10.299	1504	1511	1528	rVB	2577013	3450867	100.00%	9.540%
54	10.640	1562	1567	1580	rVB	1501741	1917616	55.57%	5.301%
55	10.963	1615	1620	1628	rVB	207113	270492	7.84%	0.748%
56	11.079	1634	1639	1645	rBV	406264	514373	14.91%	1.422%
57	11.305	1667	1676	1683	rBV	362766	441226	12.79%	1.220%
58	11.378	1683	1688	1696	rVV	1289373	2183429	63.27%	6.036%
59	11.451	1696	1700	1707	rVB	605058	712453	20.65%	1.970%
60	11.610	1721	1726	1736	rBV	867002	1101728	31.93%	3.046%
61	11.750	1744	1749	1756	rVB	2350727	2828893	81.98%	7.821%
62	11.866	1762	1768	1770	rBV	38202	54223	1.57%	0.150%
63	11.890	1770	1772	1779	rVB	57719	66445	1.93%	0.184%
64	11.969	1780	1785	1788	rBV	113179	131117	3.80%	0.362%
65	12.018	1788	1793	1799	rVV	596860	796083	23.07%	2.201%
66	12.085	1799	1804	1810	rVB	801497	937407	27.16%	2.592%
67	12.244	1824	1830	1832	rBV2	287173	381561	11.06%	1.055%
68	12.268	1832	1834	1838	rVV	327180	346239	10.03%	0.957%
69	12.317	1838	1842	1850	rVB3	379223	587593	17.03%	1.624%
70	12.402	1851	1856	1861	rVB2	48402	66910	1.94%	0.185%
71	12.463	1861	1866	1872	rVB	137628	160444	4.65%	0.444%
72	12.530	1872	1877	1879	rBV	287659	355410	10.30%	0.983%
73	12.555	1879	1881	1885	rVB	182306	189512	5.49%	0.524%
74	12.609	1885	1890	1894	rBV	580469	716418	20.76%	1.981%
75	12.646	1894	1896	1900	rVB	74279	77286	2.24%	0.214%
76	12.701	1900	1905	1912	rBV2	443831	653242	18.93%	1.806%

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ClientSampleId :
MW-3

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
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Title : SW846 8260

77	12.798	1917	1921	1925	rBV2	27126	34612	1.00%	0.096%
78	12.847	1925	1929	1934	rVB2	156129	198392	5.75%	0.548%
79	12.920	1934	1941	1945	rBV	346584	423713	12.28%	1.171%
80	12.963	1945	1948	1953	rVB	268727	306058	8.87%	0.846%
81	13.042	1953	1961	1968	rBV5	49752	122777	3.56%	0.339%
82	13.134	1972	1976	1978	rBV	37406	45191	1.31%	0.125%
83	13.170	1978	1982	1992	rVB	274938	387096	11.22%	1.070%
84	13.256	1992	1996	1997	rBV	35145	41569	1.20%	0.115%
85	13.292	1997	2002	2008	rVB2	537954	743858	21.56%	2.056%
86	13.420	2019	2023	2029	rVB	88227	117421	3.40%	0.325%
87	13.506	2032	2037	2040	rBV2	56661	80044	2.32%	0.221%
88	13.542	2040	2043	2046	rVV	134350	167336	4.85%	0.463%
89	13.585	2046	2050	2055	rVV4	125509	218653	6.34%	0.604%
90	13.640	2055	2059	2060	rVV	75495	96438	2.79%	0.267%
91	13.664	2060	2063	2068	rVB2	136959	198262	5.75%	0.548%
92	13.774	2074	2081	2088	rBV	583603	727737	21.09%	2.012%
93	13.865	2088	2096	2100	rVV3	30254	55621	1.61%	0.154%
94	13.926	2100	2106	2110	rVV2	40269	54536	1.58%	0.151%
95	13.969	2110	2113	2117	rVB	41365	48045	1.39%	0.133%
96	14.073	2126	2130	2137	rBV	62719	79958	2.32%	0.221%
97	14.213	2147	2153	2162	rBV2	54564	97944	2.84%	0.271%
98	14.371	2175	2179	2183	rVB	37283	44402	1.29%	0.123%
99	14.633	2218	2222	2233	rVB	174803	234936	6.81%	0.649%
100	14.780	2241	2246	2255	rBV	101174	137135	3.97%	0.379%

Sum of corrected areas: 36172110

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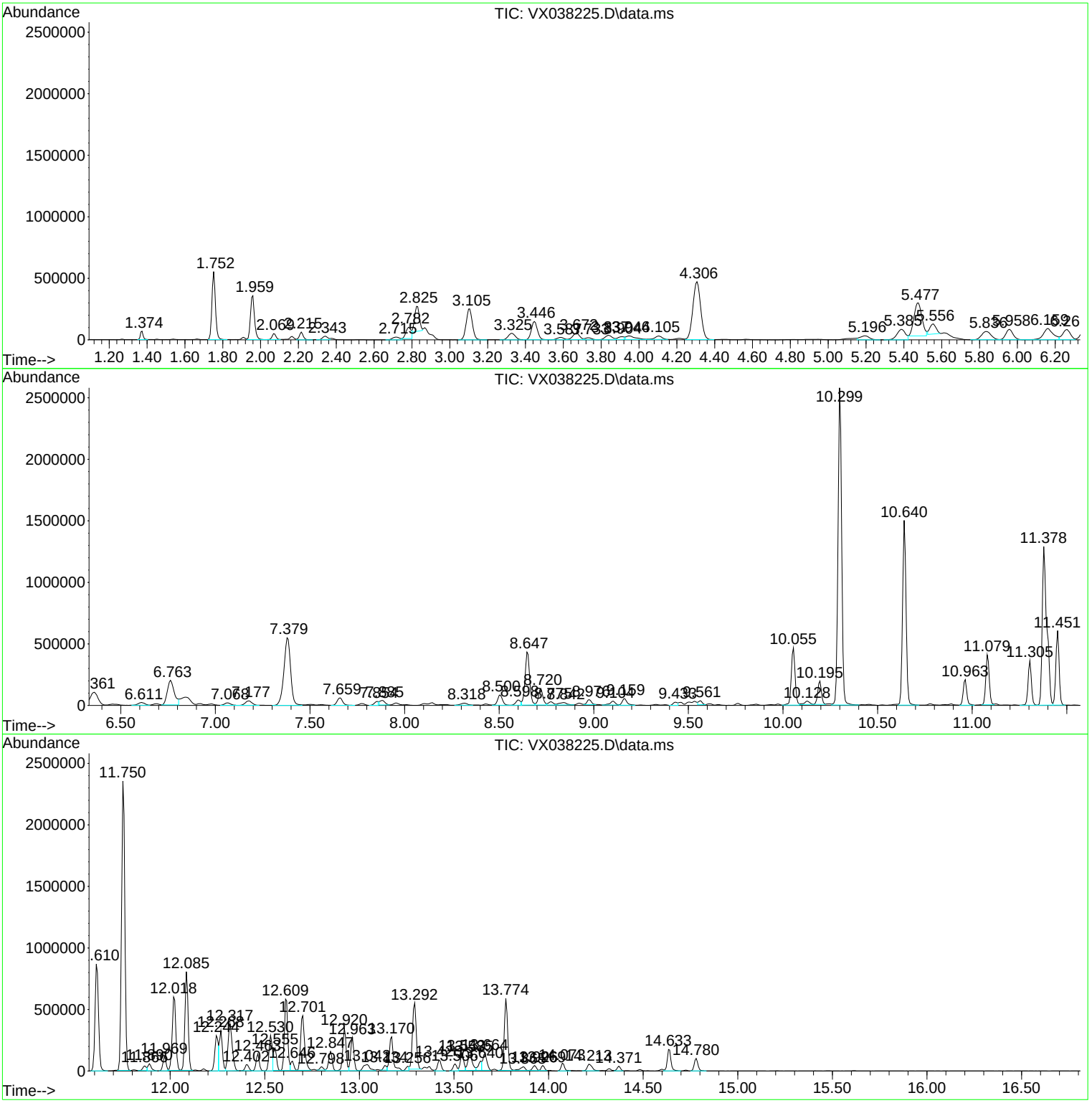
Instrument :
MSVOA_X
ClientSampleId :
MW-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.M
Quant Title : SW846 8260

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



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Instrument :
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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.M
Quant Title : SW846 8260

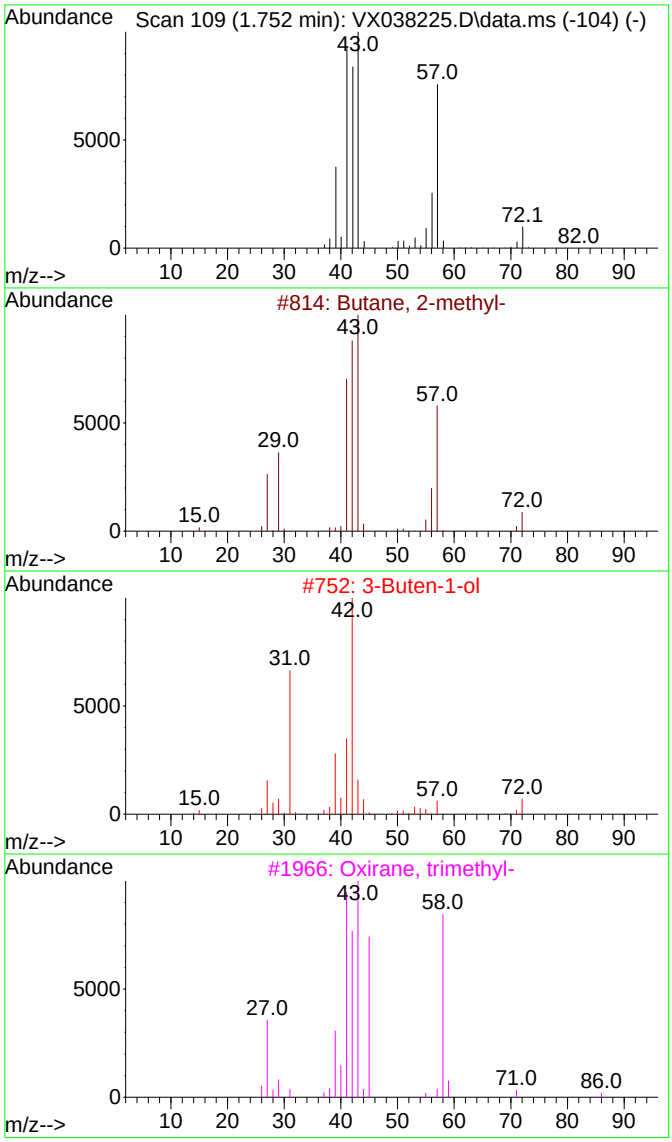
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

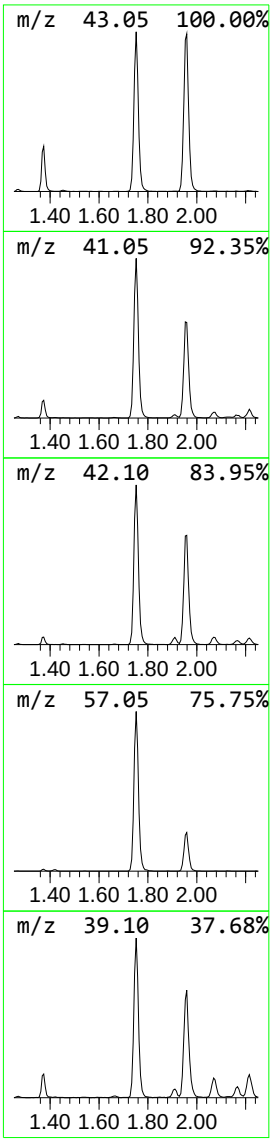
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.752	192.18 ug/l	714962	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	47
3	Oxirane, trimethyl-	86	C5H10O	005076-19-7	40
4	Pentane	72	C5H12	000109-66-0	28
5	Butane, 1-chloro-2-methyl-	106	C5H11Cl	000616-13-7	12





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101123\
Data File : VX038225.D
Acq On : 11 Oct 2023 13:36
Operator : JC/MD
Sample : 04788-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

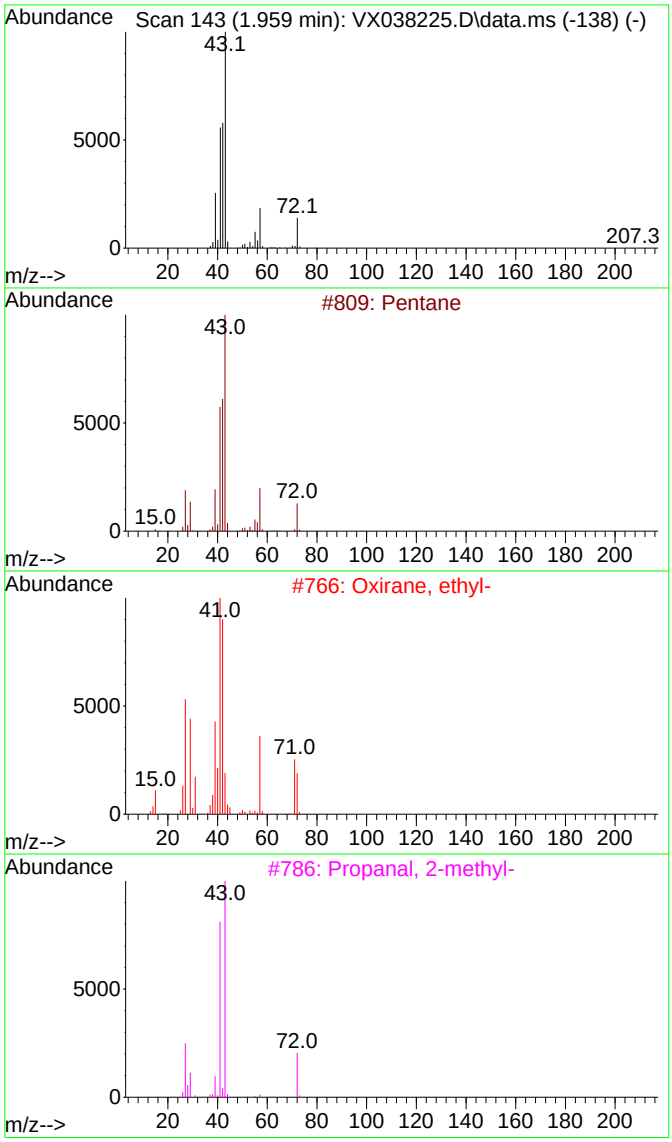
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.M
Quant Title : SW846 8260

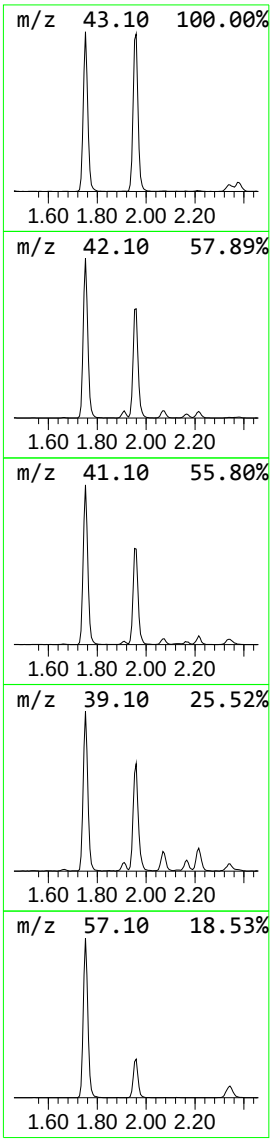
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Pentane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.959	140.10 ug/l	521198	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane	72	C5H12	000109-66-0	91
2	Oxirane, ethyl-	72	C4H8O	000106-88-7	50
3	Propanal, 2-methyl-	72	C4H8O	000078-84-2	16
4	2-Propanone, hydrazone	72	C3H8N2	005281-20-9	10
5	Diaziridine,3,3-dimethyl-	72	C3H8N2	004901-76-2	9





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Instrument :
MSVOA_X
ClientSampleId :
MW-3

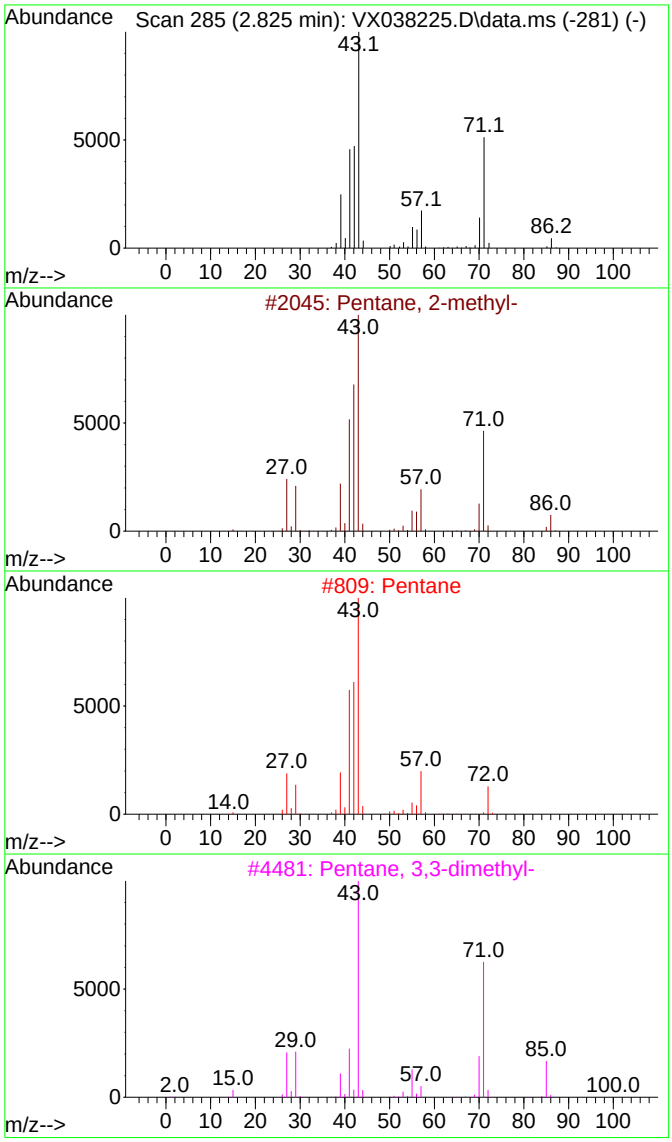
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.M
Quant Title : SW846 8260

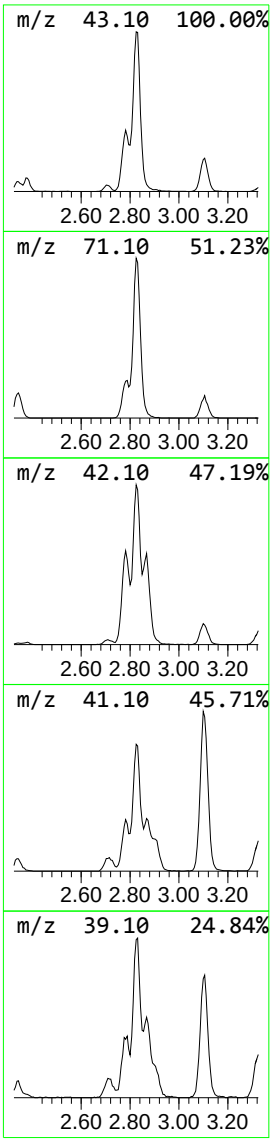
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Pentane, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.825	94.75 ug/l	352486	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	47
3	Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	38
4	1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	36
5	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	33





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ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

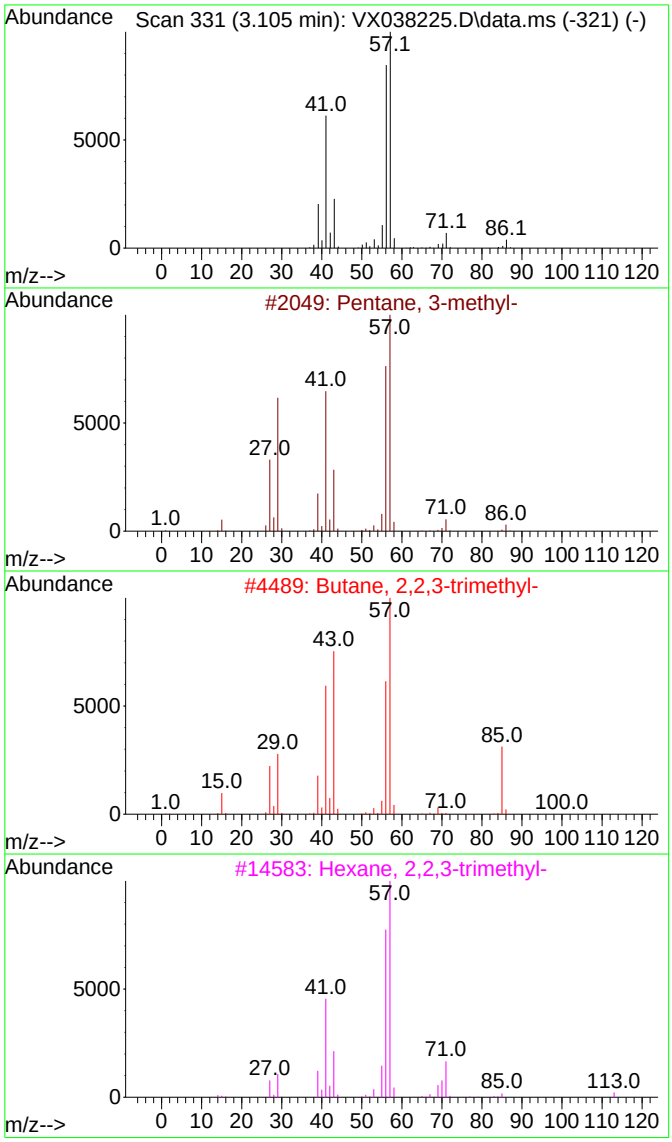
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.M
Quant Title : SW846 8260

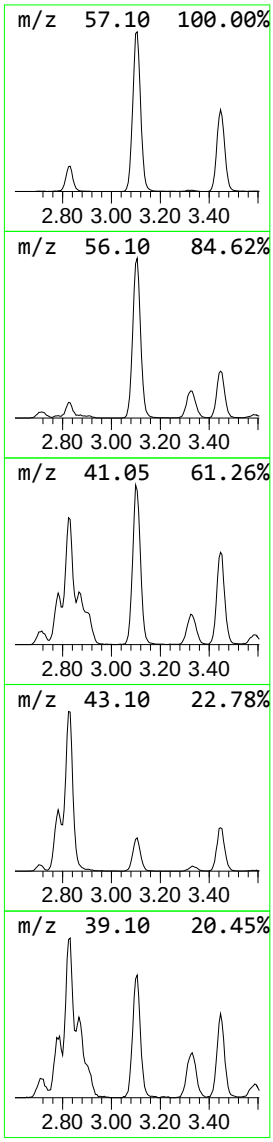
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Pentane, 3-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.105	153.12 ug/l	569621	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	90
2	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	78
3	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	78
4	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	59
5	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	43





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Data File : VX038225.D
Acq On : 11 Oct 2023 13:36
Operator : JC/MD
Sample : 04788-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.M
Quant Title : SW846 8260

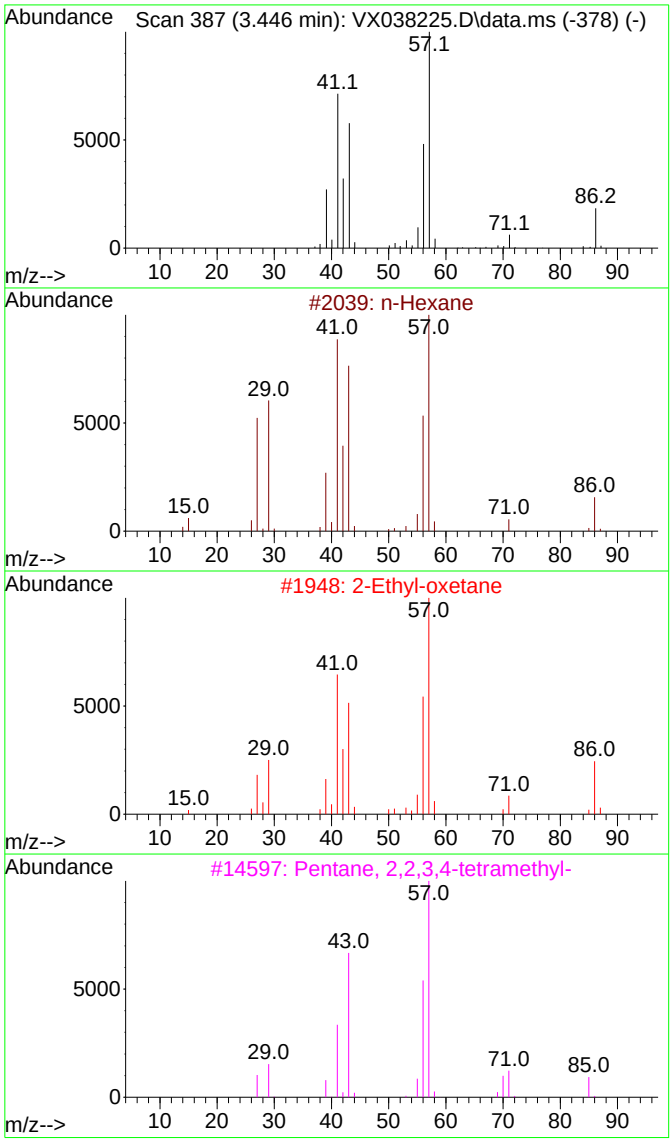
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

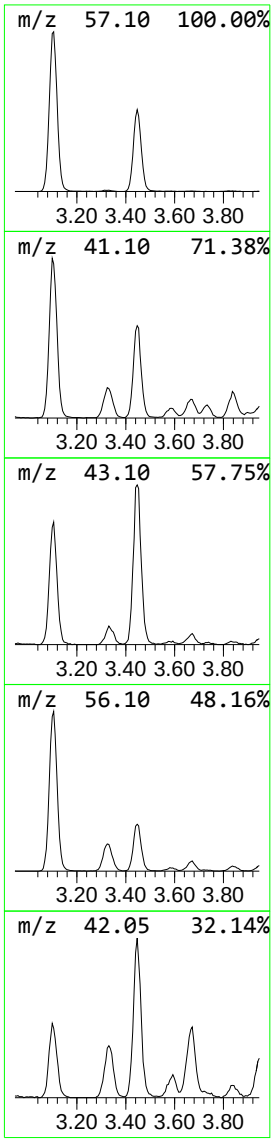
****

Peak Number 5 n-Hexane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.446	89.40 ug/l	332603	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexane	86	C6H14	000110-54-3	91
2	2-Ethyl-oxetane	86	C5H10O	1010386-40-2	90
3	Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	50
4	1-Pentene, 4-methyl-	84	C6H12	000691-37-2	38
5	Butanal, 2-methyl-	86	C5H10O	000096-17-3	35





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101123\
Data File : VX038225.D
Acq On : 11 Oct 2023 13:36
Operator : JC/MD
Sample : 04788-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

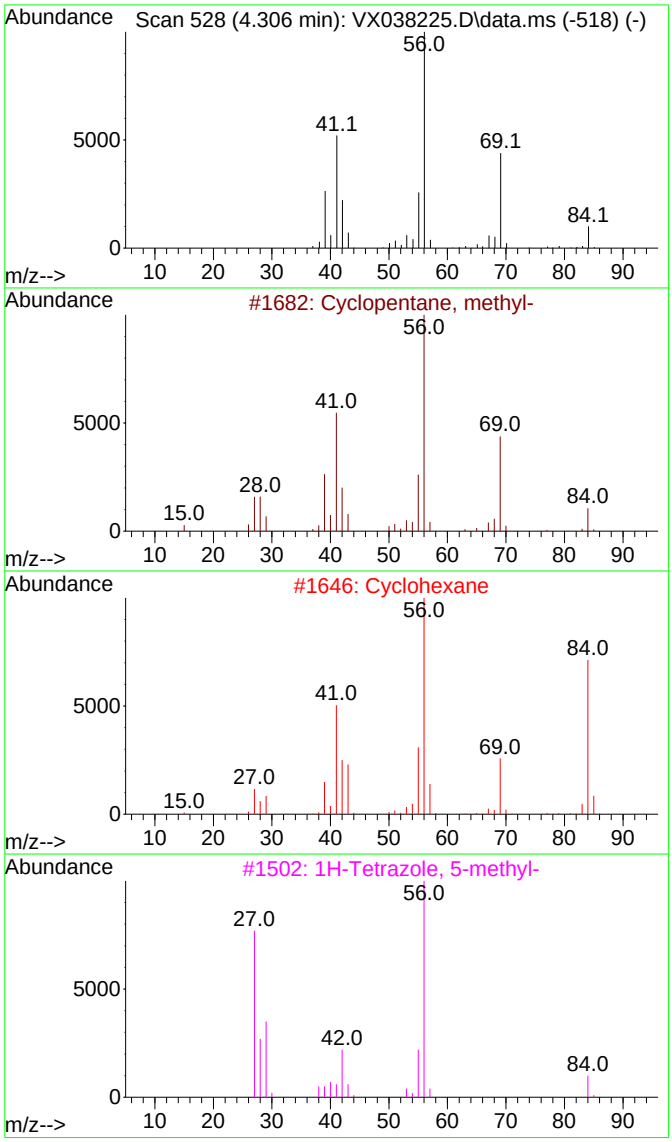
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.M
Quant Title : SW846 8260

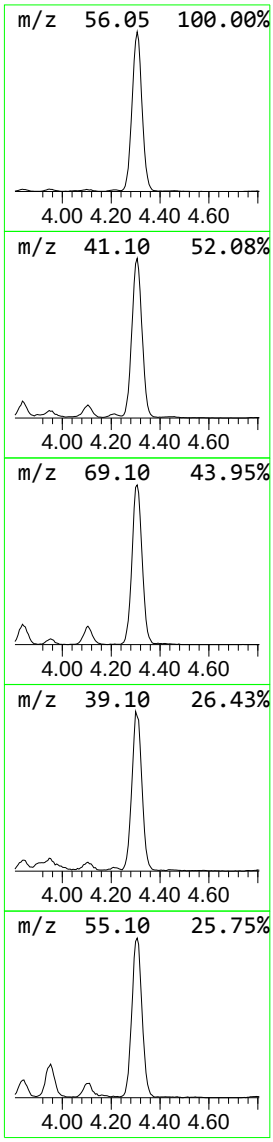
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Cyclopentane, methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.306	360.56 ug/l	1341380	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2	Cyclohexane	84	C6H12	000110-82-7	78
3	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
4	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
5	Cyclobutanecarbonitrile, 3,3-dim...	109	C7H11N	053783-86-1	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101123\
Data File : VX038225.D
Acq On : 11 Oct 2023 13:36
Operator : JC/MD
Sample : 04788-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

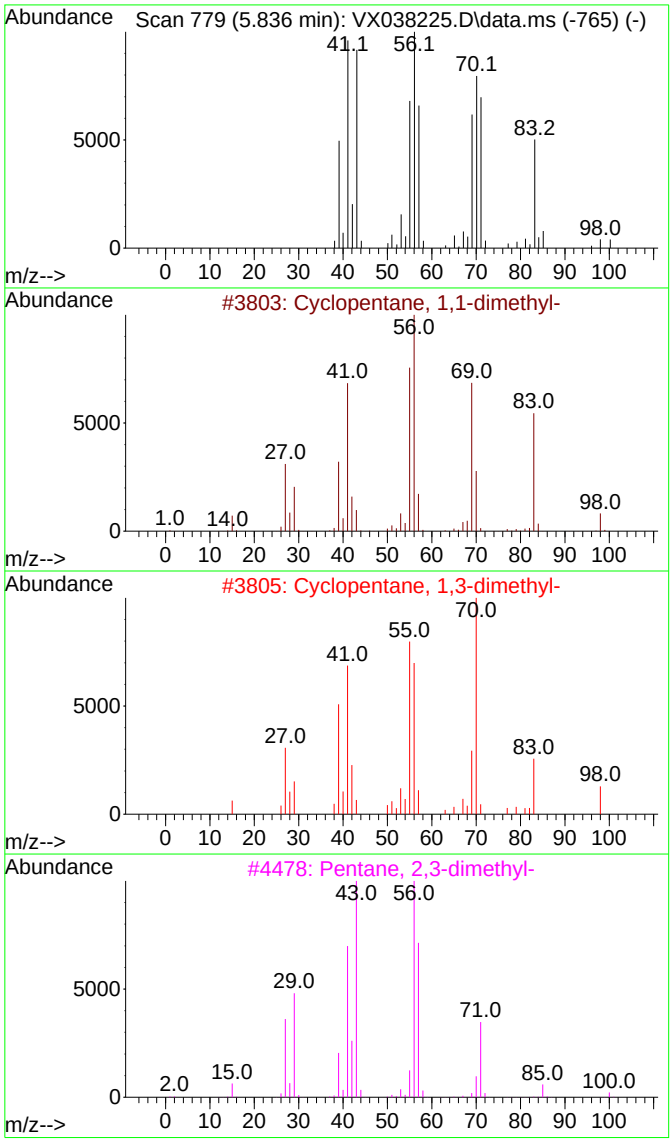
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.M
Quant Title : SW846 8260

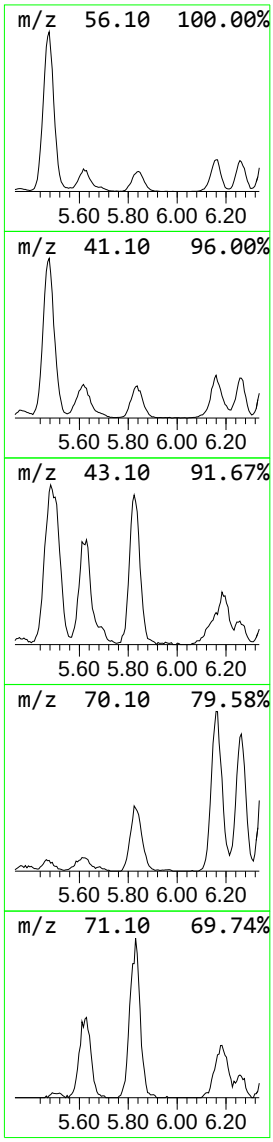
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 unknown5.836 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.836	63.41 ug/l	235917	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	45
2	Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	38
3	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	38
4	2-Heptene	98	C7H14	000592-77-8	38
5	1-Heptene, 4-methyl-	112	C8H16	013151-05-8	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101123\
Data File : VX038225.D
Acq On : 11 Oct 2023 13:36
Operator : JC/MD
Sample : 04788-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

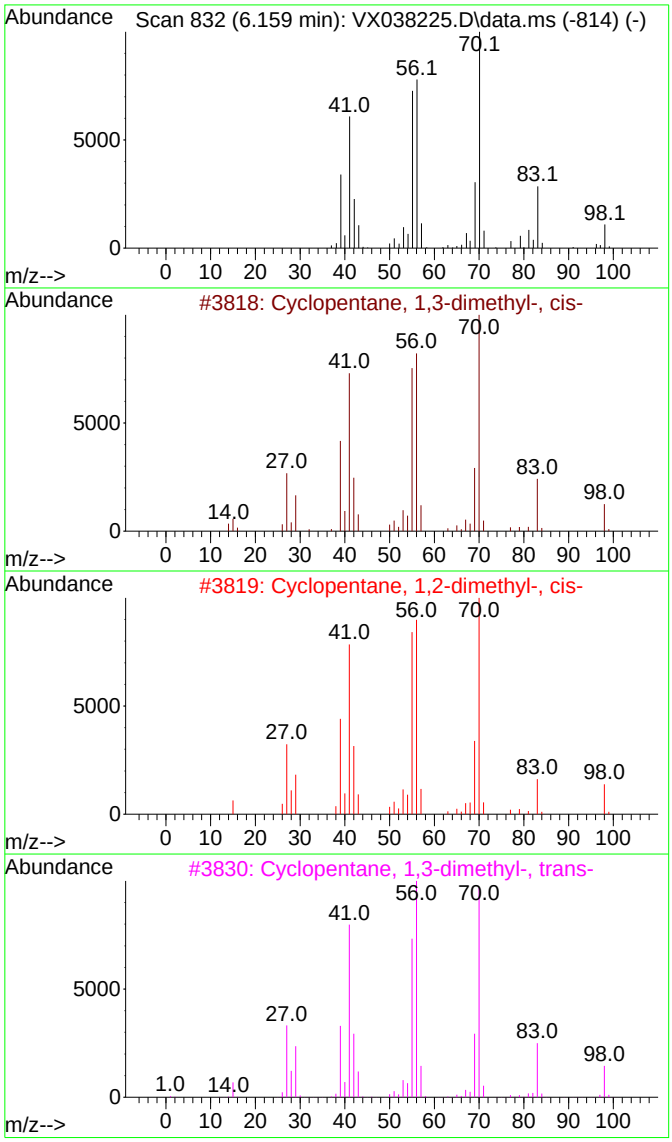
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.M
Quant Title : SW846 8260

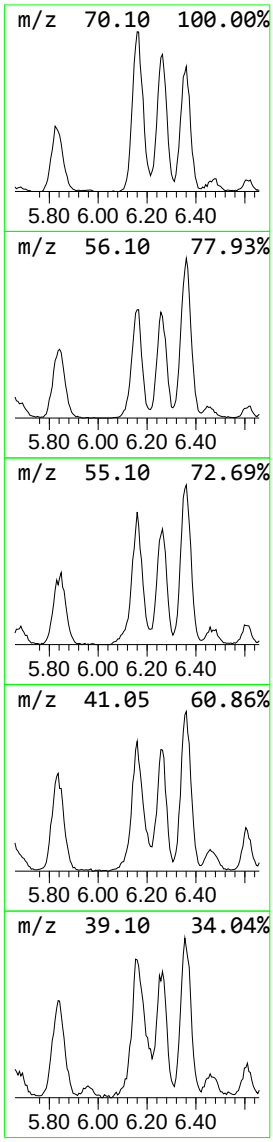
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Cyclopentane, 1,3-dimethyl-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.159	89.65 ug/l	333507	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	96
2	Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	94
3	Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	94
4	Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	93
5	Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101123\
Data File : VX038225.D
Acq On : 11 Oct 2023 13:36
Operator : JC/MD
Sample : 04788-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

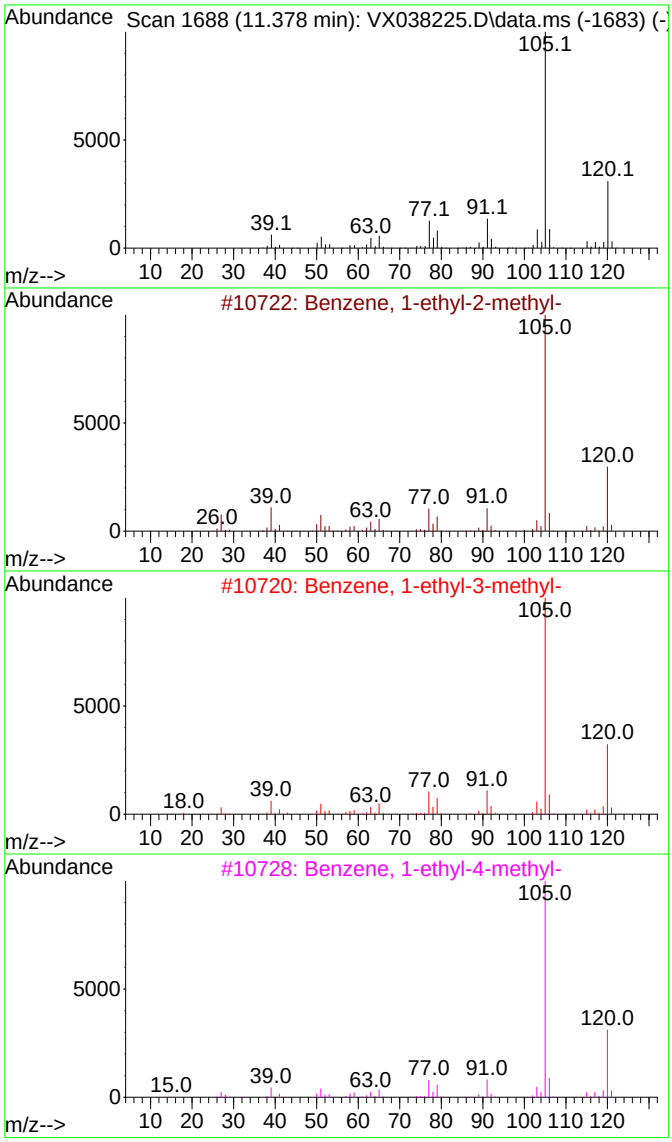
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.M
Quant Title : SW846 8260

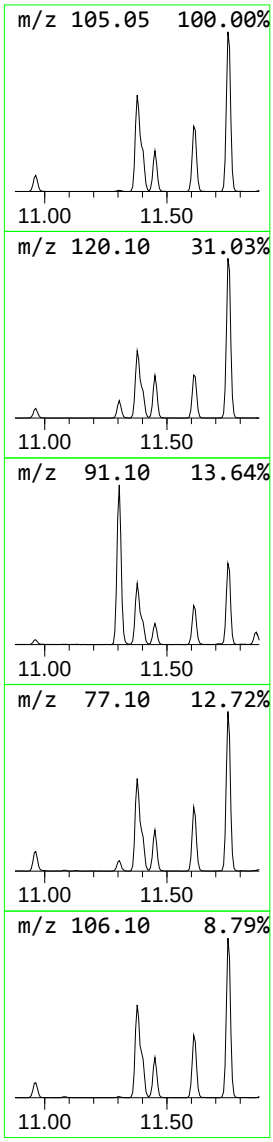
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1-ethyl-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.378	137.14 ug/l	2183430	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	94
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Mesitylene	120	C9H12	000108-67-8	91





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Data File : VX038225.D
Acq On : 11 Oct 2023 13:36
Operator : JC/MD
Sample : 04788-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

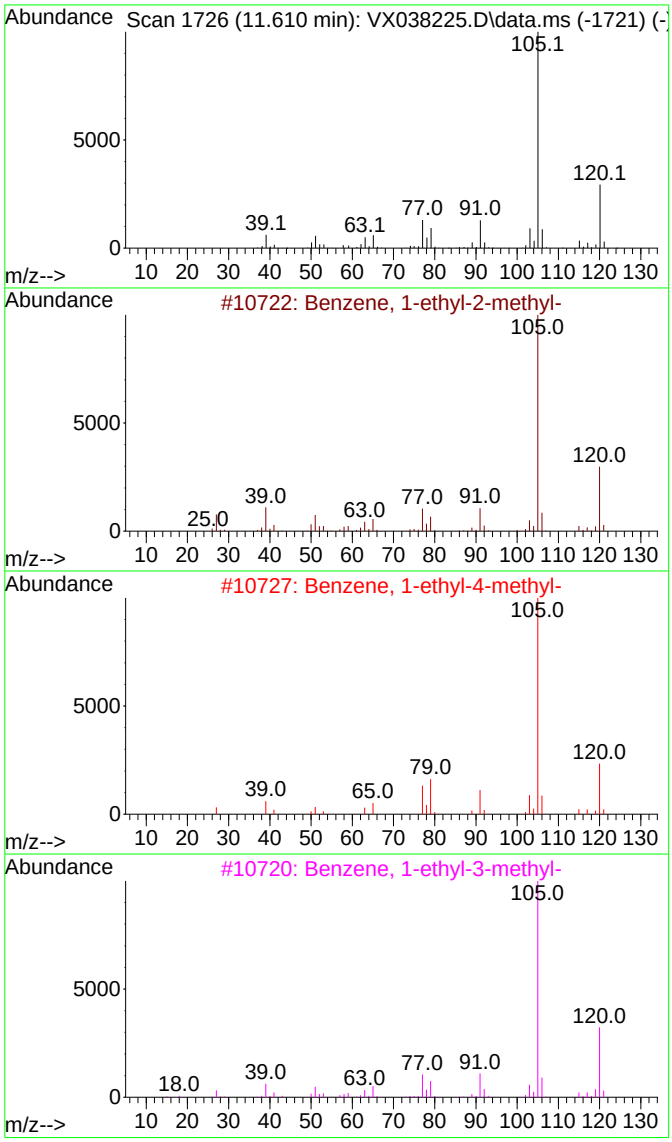
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.M
Quant Title : SW846 8260

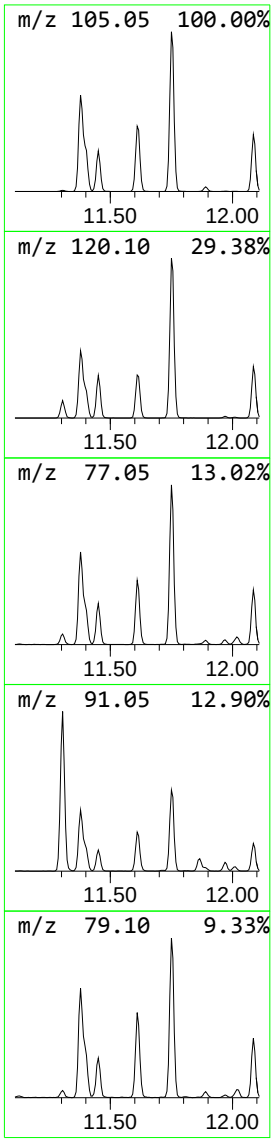
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 1-ethyl-4-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.610	69.20 ug/l	1101730	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	94
2	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Mesitylene	120	C9H12	000108-67-8	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101123\
Data File : VX038225.D
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ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.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-methyl-	1.752	192.2	ug/l	714962	1	5.556	186011	50.0
Pentane	1.959	140.1	ug/l	521198	1	5.556	186011	50.0
Pentane, 2-methyl-	2.825	94.8	ug/l	352486	1	5.556	186011	50.0
Pentane, 3-methyl-	3.105	153.1	ug/l	569621	1	5.556	186011	50.0
n-Hexane	3.446	89.4	ug/l	332603	1	5.556	186011	50.0
Cyclopentane, m...	4.306	360.6	ug/l	1341380	1	5.556	186011	50.0
unknown5.836	5.836	63.4	ug/l	235917	1	5.556	186011	50.0
Cyclopentane, 1...	6.159	89.7	ug/l	333507	1	5.556	186011	50.0
Benzene, 1-ethy...	11.378	137.1	ug/l	2183430	4	12.024	796083	50.0
Benzene, 1-ethy...	11.610	69.2	ug/l	1101730	4	12.024	796083	50.0