LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
Data File : VX028662.D
Acq On : 13 May 2022 00:39
Operator : JC/MD
Sample : N2811-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 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\82X051222W.M
Title : SW846 8260

Signal : TIC: VX028662.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.752	103	109	118	rBV	362364	501295	7.85%	1.090%
2	1.953	137	142	155	rVB	533134	749904	11.75%	1.630%
3	2.069	156	161	170	rVV	52176	72150	1.13%	0.157%
4	2.709	256	266	271	rBV4	34668	83903	1.31%	0.182%
5	2.782	272	278	280	rBV3	47303	83766	1.31%	0.182%
6	2.825	280	285	290	rBV	269553	505819	7.92%	1.100%
7	3.105	321	331	343	rBV	222576	500965	7.85%	1.089%
8	3.447	377	387	398	rBV	357751	798659	12.51%	1.736%
9	3.672	416	424	429	rVV	38653	92851	1.45%	0.202%
10	3.733	429	434	445	rVB3	30524	73468	1.15%	0.160%
11	4.306	518	528	544	rVB	474424	1360868	21.32%	2.958%
12	5.391	690	706	712	rBV	165931	496403	7.78%	1.079%
13	5.477	712	720	728	rVV3	378001	1215714	19.04%	2.643%
14	5.556	728	733	740	rVB	192114	480754	7.53%	1.045%
15	5.836	765	779	791	rBV4	154186	522854	8.19%	1.137%
16	5.958	791	799	804	rVV2	174346	446238	6.99%	0.970%
17	6.044	804	813	824	rVV	2427409	6384200	100.00%	13.878%
18	6.159	824	832	842	rVV3	159001	549499	8.61%	1.195%
19	6.263	842	849	857	rVV	145079	407766	6.39%	0.886%
20	6.361	857	865	879	rVB2	235399	659612	10.33%	1.434%
21	6.617	895	907	915	rBV	162855	399661	6.26%	0.869%
22	6.763	923	931	941	rBV	422297	1019696	15.97%	2.217%
23	7.385	1019	1033	1045	rBV	1207288	2842226	44.52%	6.178%
24	7.659	1071	1078	1089	rVB	149858	295051	4.62%	0.641%
25	7.775	1089	1097	1104	rBV	74475	151162	2.37%	0.329%
26	7.958	1121	1127	1141	rVB2	100831	220390	3.45%	0.479%
27	8.275	1175	1179	1183	rBV	65254	109649	1.72%	0.238%
28	8.318	1183	1186	1193	rVB3	59194	104616	1.64%	0.227%
29	8.440	1201	1206	1210	rBV	44000	76101	1.19%	0.165%
30	8.507	1212	1217	1226	rVB2	77514	155198	2.43%	0.337%
31	8.604	1226	1233	1236	rBV2	249568	473853	7.42%	1.030%
32	8.653	1236	1241	1247	rVV	916624	1549242	24.27%	3.368%
33	8.720	1247	1252	1257	rVV	208492	328829	5.15%	0.715%
34	8.775	1257	1261	1265	rVV	119794	196290	3.07%	0.427%
35	8.824	1265	1269	1270	rVV2	44867	66525	1.04%	0.145%
36	8.848	1270	1273	1279	rVV2	85936	137555	2.15%	0.299%

LSC Area Percent Report

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Operator : JC/MD
Sample : N2811-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 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\82X051222W.M
Title : SW846 8260

37	8.915	1279	1284	1289	rVV2	98245	151511	2.37%	0.329%
38	8.976	1289	1294	1305	rVB	211034	373864	5.86%	0.813%
39	9.104	1305	1315	1320	rBV	111487	199684	3.13%	0.434%
40	9.159	1320	1324	1329	rVV2	45727	76029	1.19%	0.165%
41	9.512	1377	1382	1386	rBV4	100394	163488	2.56%	0.355%
42	9.567	1386	1391	1395	rVB	246981	348396	5.46%	0.757%
43	9.622	1395	1400	1404	rBV	156888	238327	3.73%	0.518%
44	9.762	1418	1423	1428	rVB2	41092	68130	1.07%	0.148%
45	9.836	1428	1435	1443	rBV5	48677	121629	1.91%	0.264%
46	10.055	1466	1471	1476	rVV	960469	1240735	19.43%	2.697%
47	10.128	1476	1483	1488	rVB	90392	155210	2.43%	0.337%
48	10.195	1488	1494	1499	rBV	476569	626102	9.81%	1.361%
49	10.305	1502	1512	1518	rVV	1841485	2694893	42.21%	5.858%
50	10.360	1518	1521	1532	rVB3	62417	102033	1.60%	0.222%
51	10.592	1551	1559	1565	rBV4	67182	161170	2.52%	0.350%
52	10.646	1565	1568	1578	rVB2	45312	78099	1.22%	0.170%
53	10.781	1578	1590	1595	rBV2	133260	218186	3.42%	0.474%
54	10.860	1597	1603	1615	rVB4	88253	158788	2.49%	0.345%
55	10.963	1615	1620	1624	rBV	200844	254258	3.98%	0.553%
56	11.085	1634	1640	1644	rBV	762945	1003813	15.72%	2.182%
57	11.128	1644	1647	1651	rVB	51027	66493	1.04%	0.145%
58	11.226	1659	1663	1668	rVB4	46527	68135	1.07%	0.148%
59	11.305	1669	1676	1683	rBV	246026	310023	4.86%	0.674%
60	11.378	1683	1688	1696	rVV	696618	1289175	20.19%	2.802%
61	11.451	1696	1700	1709	rVB	546827	701089	10.98%	1.524%
62	11.616	1722	1727	1735	rVB	253279	351655	5.51%	0.764%
63	11.750	1745	1749	1755	rVB	1461103	1833009	28.71%	3.985%
64	11.866	1761	1768	1770	rBV	49682	77329	1.21%	0.168%
65	11.890	1770	1772	1777	rVB	89142	104523	1.64%	0.227%
66	11.969	1777	1785	1789	rBV	179147	240561	3.77%	0.523%
67	12.024	1789	1794	1799	rVB	1068347	1400060	21.93%	3.043%
68	12.085	1799	1804	1810	rBV	667952	831007	13.02%	1.806%
69	12.244	1825	1830	1832	rBV3	261620	341137	5.34%	0.742%
70	12.268	1832	1834	1838	rVV	273562	290990	4.56%	0.633%
71	12.317	1838	1842	1851	rVB4	354622	546131	8.55%	1.187%
72	12.408	1851	1857	1862	rBV3	57289	89853	1.41%	0.195%
73	12.463	1862	1866	1872	rVB	164117	191403	3.00%	0.416%
74	12.530	1872	1877	1879	rBV	186339	239782	3.76%	0.521%
75	12.555	1879	1881	1885	rVB	232346	246109	3.85%	0.535%
76	12.616	1886	1891	1894	rBV	258384	306628	4.80%	0.667%

LSC Area Percent Report

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Sample : N2811-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 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

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Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Title : SW846 8260

77	12.646	1894	1896	1900	rVB	91673	97038	1.52%	0.211%
78	12.701	1900	1905	1913	rBV3	327390	540252	8.46%	1.174%
79	12.847	1924	1929	1934	rVB2	124769	174372	2.73%	0.379%
80	12.920	1934	1941	1945	rBV	126901	189900	2.97%	0.413%
81	12.963	1945	1948	1954	rVB	228006	277748	4.35%	0.604%
82	13.030	1954	1959	1964	rBV2	65153	131623	2.06%	0.286%
83	13.140	1973	1977	1979	rBV2	58091	79908	1.25%	0.174%
84	13.170	1979	1982	1985	rVV2	103398	131360	2.06%	0.286%
85	13.250	1992	1995	1998	rBV2	53328	81158	1.27%	0.176%
86	13.292	1998	2002	2007	rVB2	468710	650402	10.19%	1.414%
87	13.426	2020	2024	2031	rVB2	85664	123101	1.93%	0.268%
88	13.542	2040	2043	2046	rVV	81246	108712	1.70%	0.236%
89	13.585	2046	2050	2054	rVV3	99674	172925	2.71%	0.376%
90	13.664	2054	2063	2069	rVB3	130908	287902	4.51%	0.626%
91	13.774	2074	2081	2090	rVB	239460	350103	5.48%	0.761%
92	13.926	2099	2106	2110	rBV4	36406	73504	1.15%	0.160%
93	14.213	2148	2153	2162	rVB3	51237	91999	1.44%	0.200%
94	14.634	2210	2222	2227	rBV	169379	236492	3.70%	0.514%
95	14.780	2240	2246	2257	rVB	94089	131564	2.06%	0.286%

Sum of corrected areas: 46002232

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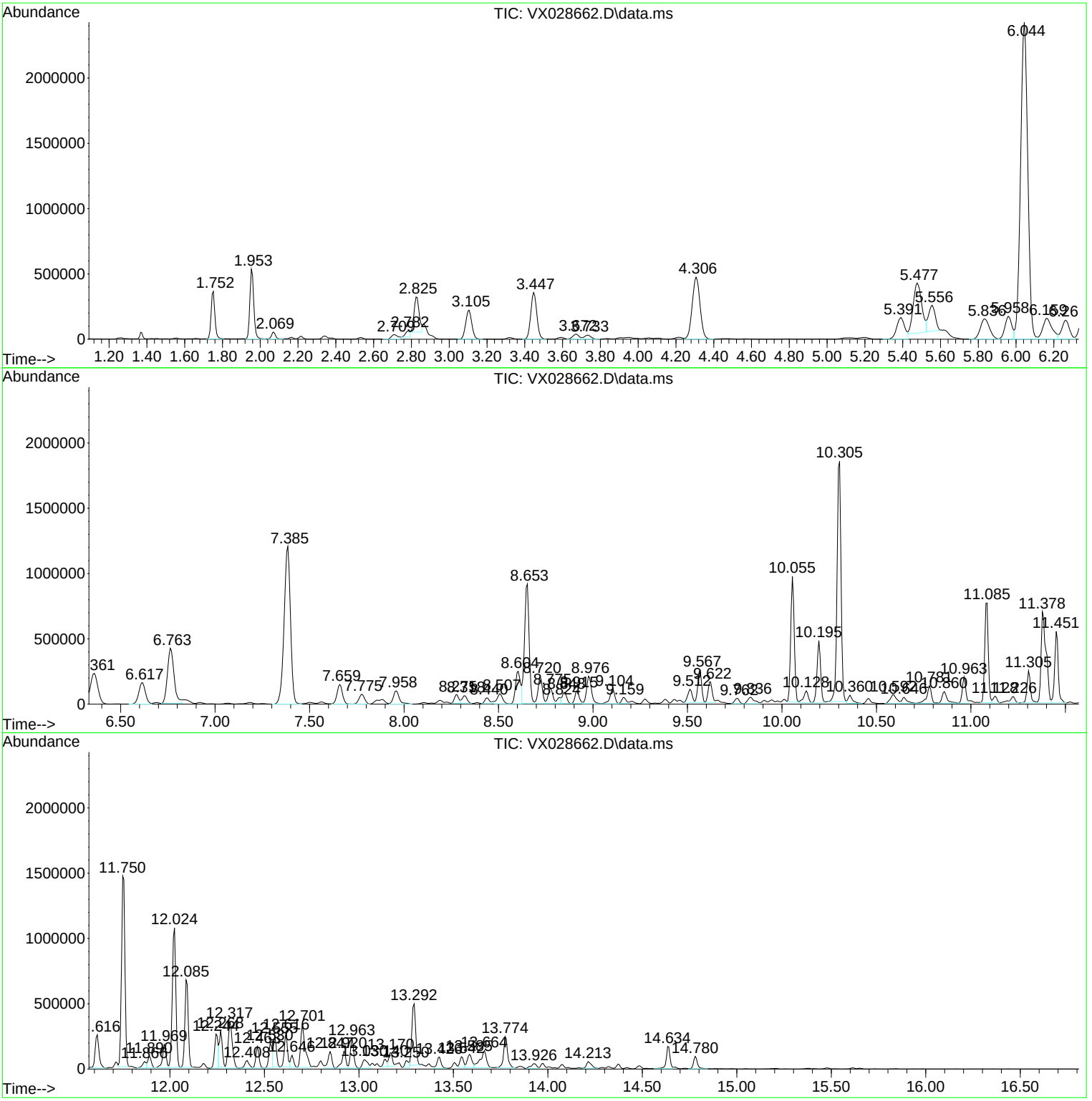
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
Data File : VX028662.D
Acq On : 13 May 2022 00:39
Operator : JC/MD
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ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

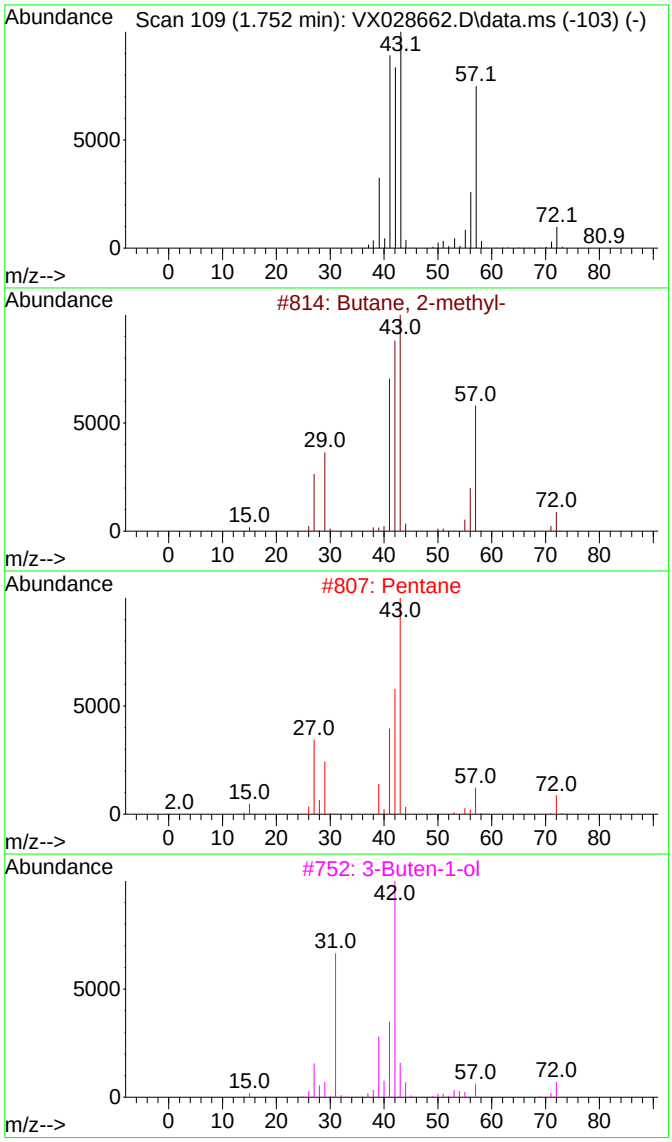
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

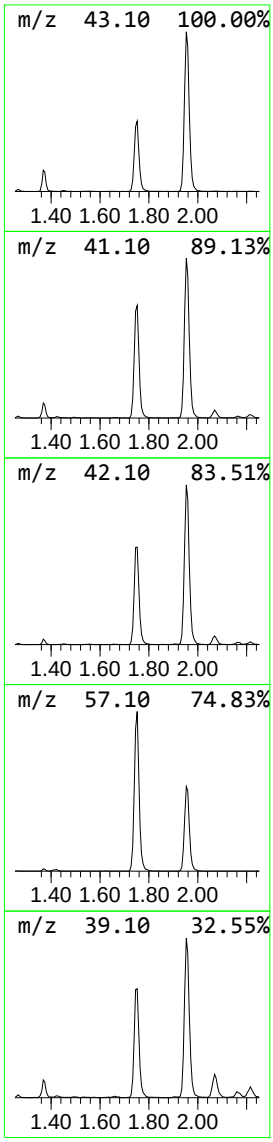
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Butane, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.752	52.14 ug/l	501295	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	91
2	Pentane	72	C5H12	000109-66-0	45
3	3-Buten-1-ol	72	C4H8O	000627-27-0	43
4	2(3H)-Furanone, dihydro-4-methyl-	100	C5H8O2	001679-49-8	28
5	1-Propene, 2-methyl-	56	C4H8	000115-11-7	10





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
Data File : VX028662.D
Acq On : 13 May 2022 00:39
Operator : JC/MD
Sample : N2811-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

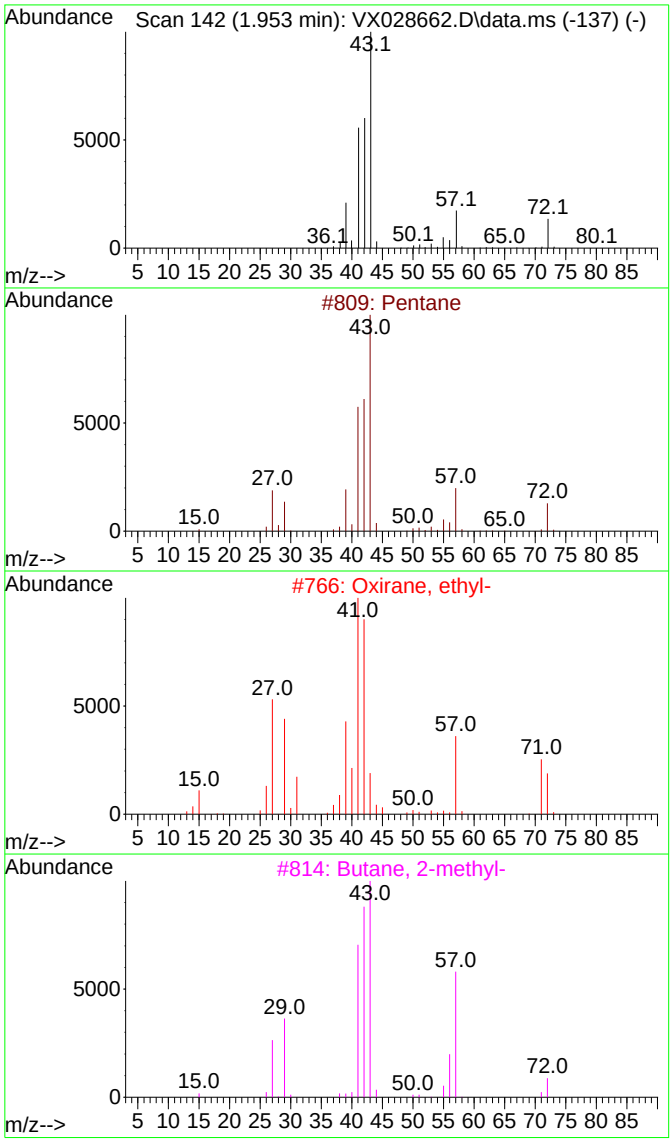
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

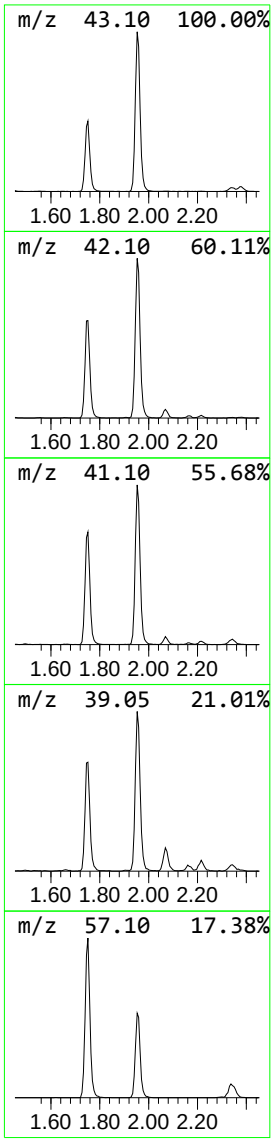
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Pentane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.953	77.99 ug/l	749904	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	Butane, 2-methyl-	72	C5H12	000078-78-4	38
4	Diaziridine,3,3-dimethyl-	72	C3H8N2	004901-76-2	35
5	2(3H)-Furanone, dihydro-4-methyl-	100	C5H8O2	001679-49-8	25





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
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Acq On : 13 May 2022 00:39
Operator : JC/MD
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

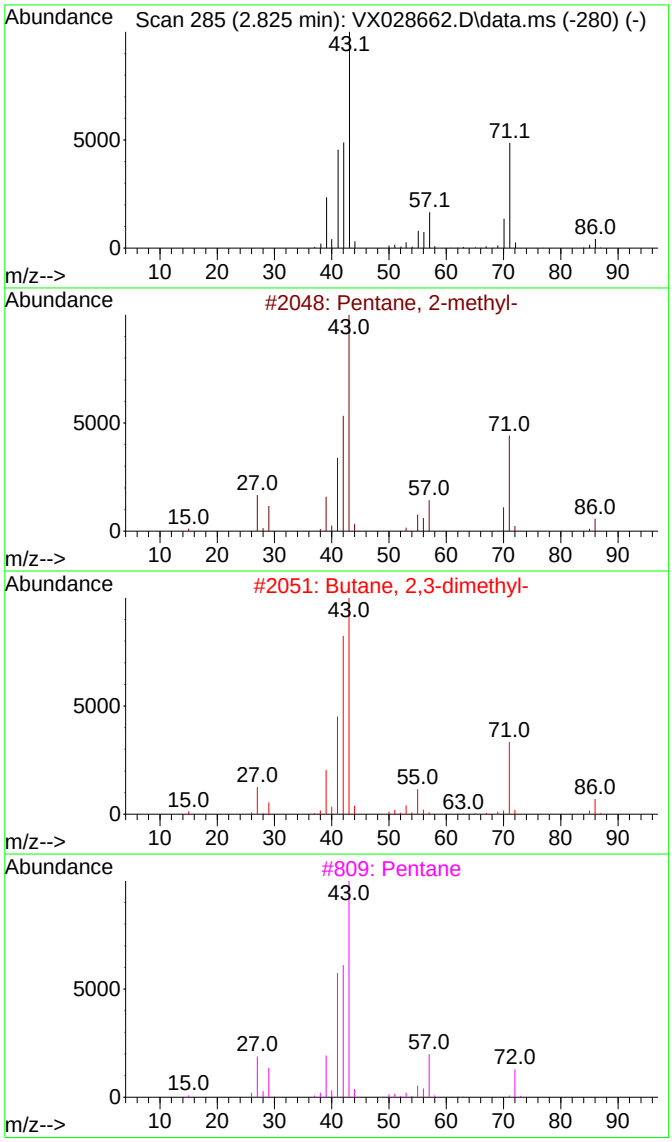
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

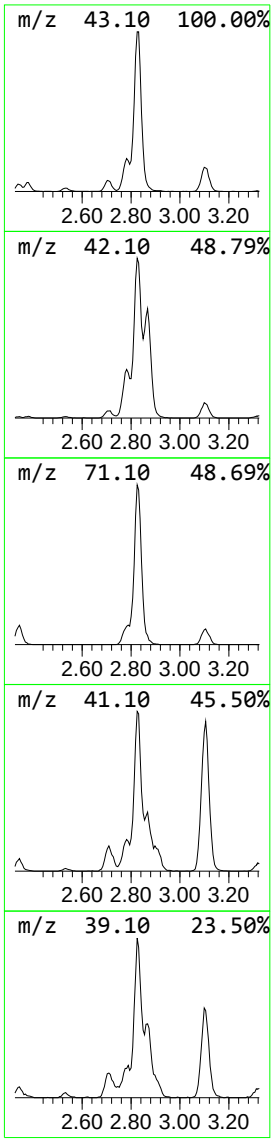
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.825	52.61 ug/l	505819	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2-methyl-	86	C6H14	000107-83-5	91
2	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	50
3	Pentane	72	C5H12	000109-66-0	46
4	1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	45
5	Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
Data File : VX028662.D
Acq On : 13 May 2022 00:39
Operator : JC/MD
Sample : N2811-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

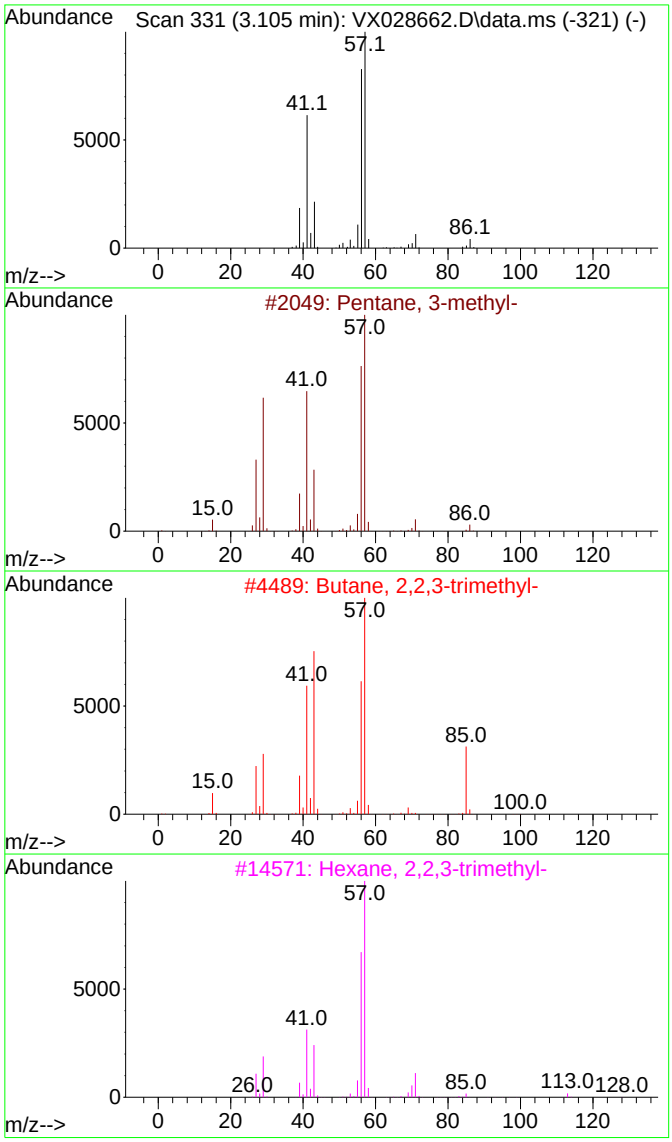
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

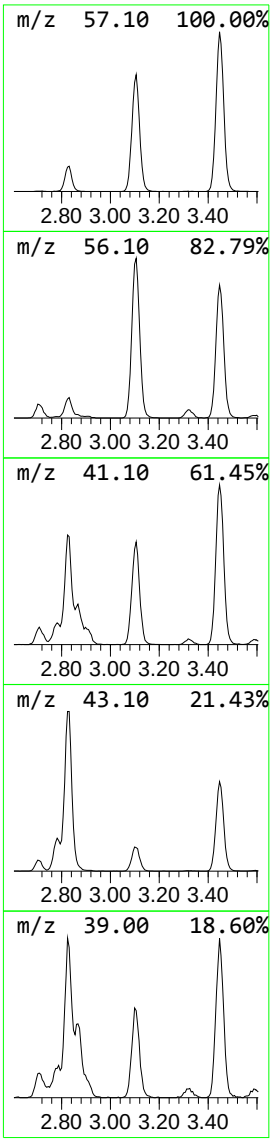
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Pentane, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.105	52.10 ug/l	500965	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
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	3,4-Dimethyldihydrofuran-2,5-dione	128	C6H8O3	007475-92-5	53
5	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	45





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
Data File : VX028662.D
Acq On : 13 May 2022 00:39
Operator : JC/MD
Sample : N2811-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

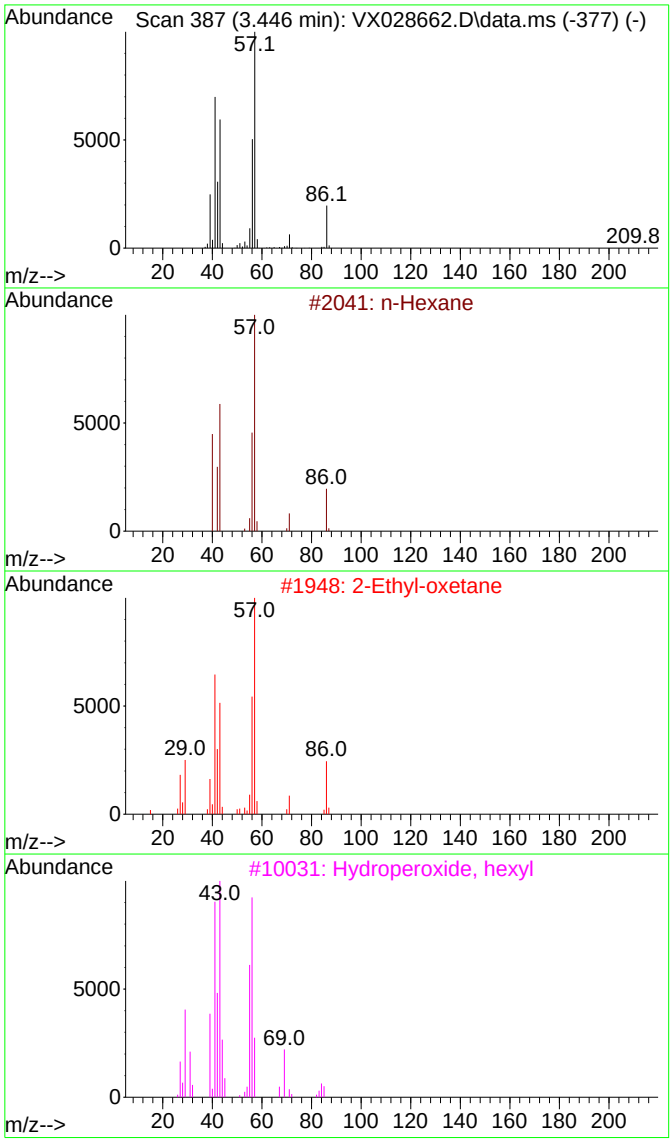
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

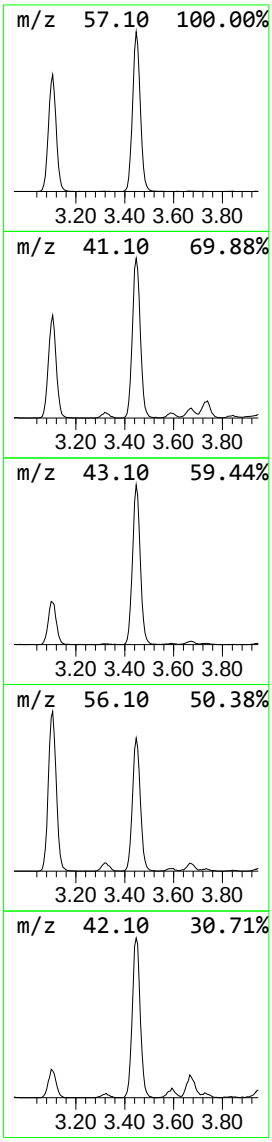
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 n-Hexane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.446	83.06 ug/l	798659	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexane	86	C6H14	000110-54-3	87
2	2-Ethyl-oxetane	86	C5H10O	1010386-40-2	86
3	Hydroperoxide, hexyl	118	C6H14O2	004312-76-9	38
4	Morpholine	87	C4H9NO	000110-91-8	38
5	1-Pentene, 4-methyl-	84	C6H12	000691-37-2	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
Data File : VX028662.D
Acq On : 13 May 2022 00:39
Operator : JC/MD
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

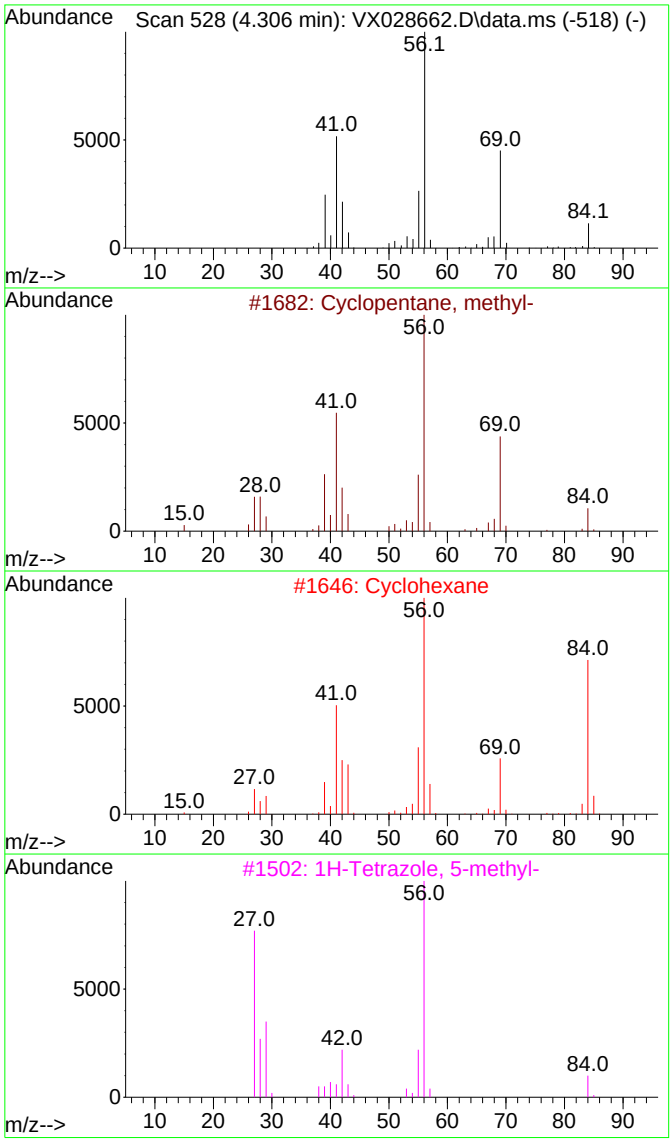
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

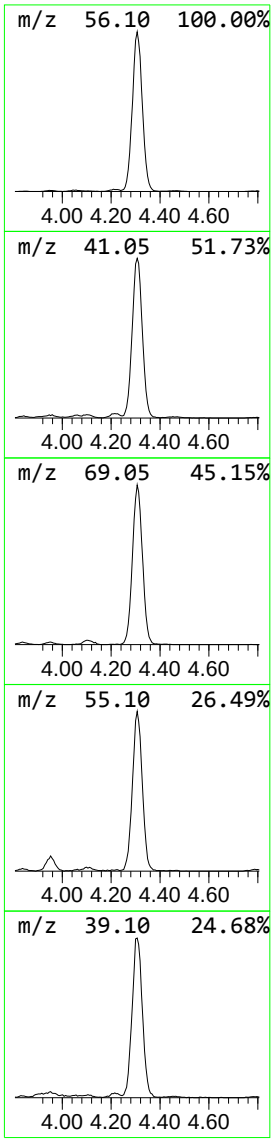
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.306	141.54 ug/l	1360870	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	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	53





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
Data File : VX028662.D
Acq On : 13 May 2022 00:39
Operator : JC/MD
Sample : N2811-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

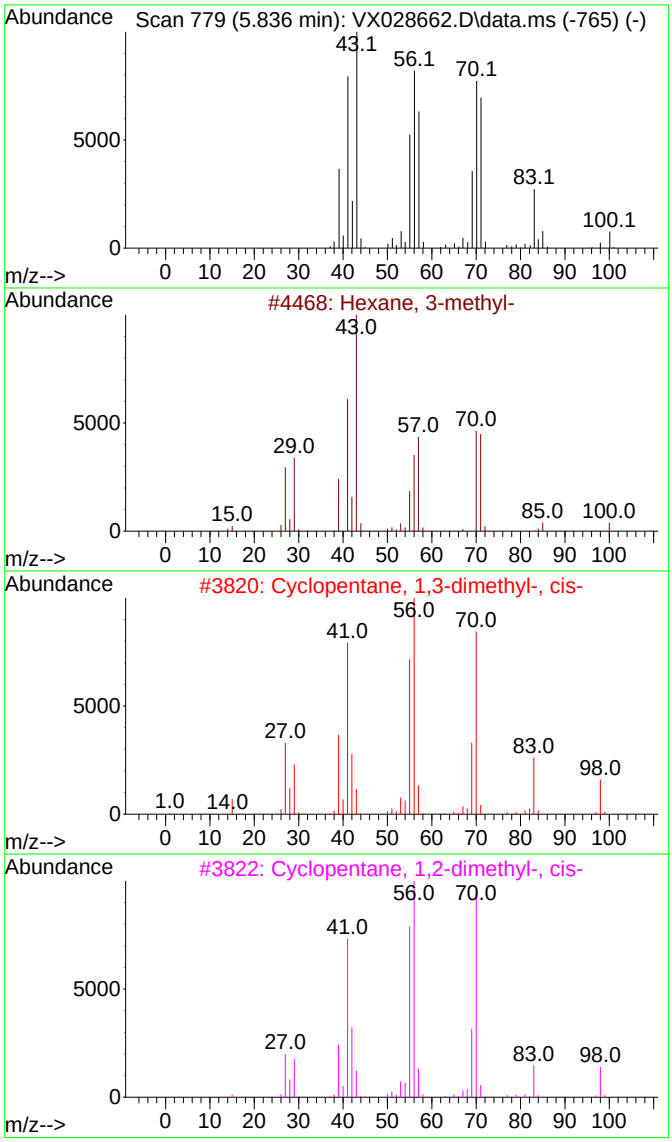
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

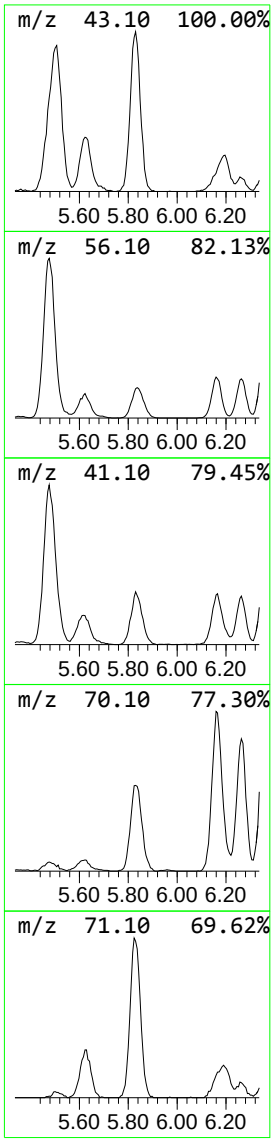
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Hexane, 3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.836	54.38 ug/l	522854	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 3-methyl-	100	C7H16	000589-34-4	70
2	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	64
3	Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	59
4	Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	47
5	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	47





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
Data File : VX028662.D
Acq On : 13 May 2022 00:39
Operator : JC/MD
Sample : N2811-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

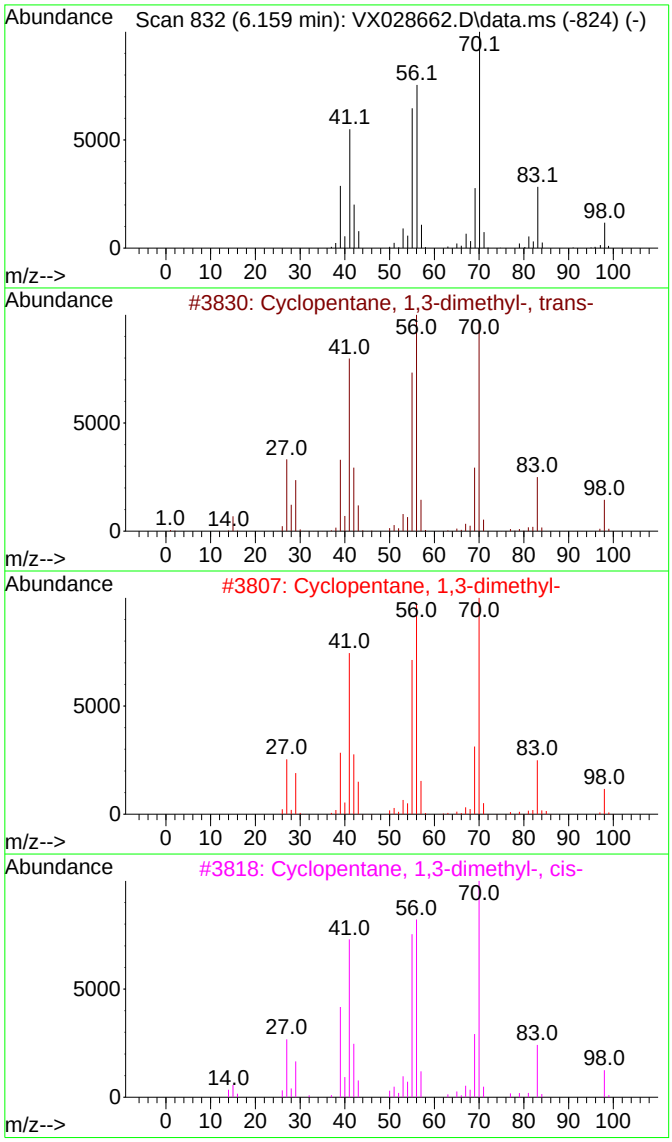
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

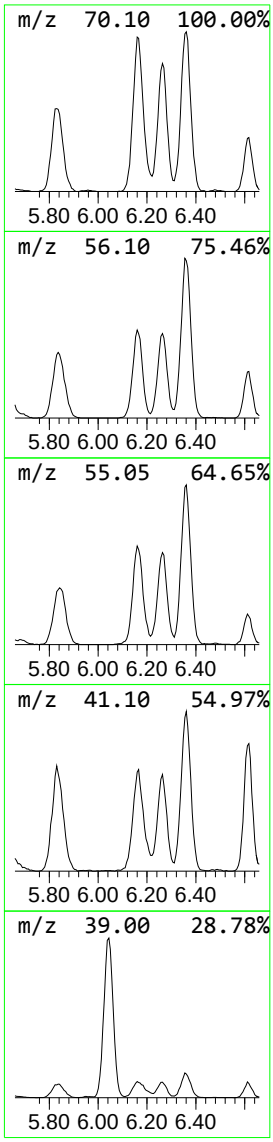
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Cyclopentane, 1,3-dimethyl-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.159	57.15 ug/l	549499	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	94
2	Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	93
3	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	93
4	Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	76
5	2H-Pyran-2,6(3H)-dione, dihydro-...	142	C7H10O3	004160-82-1	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
Data File : VX028662.D
Acq On : 13 May 2022 00:39
Operator : JC/MD
Sample : N2811-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

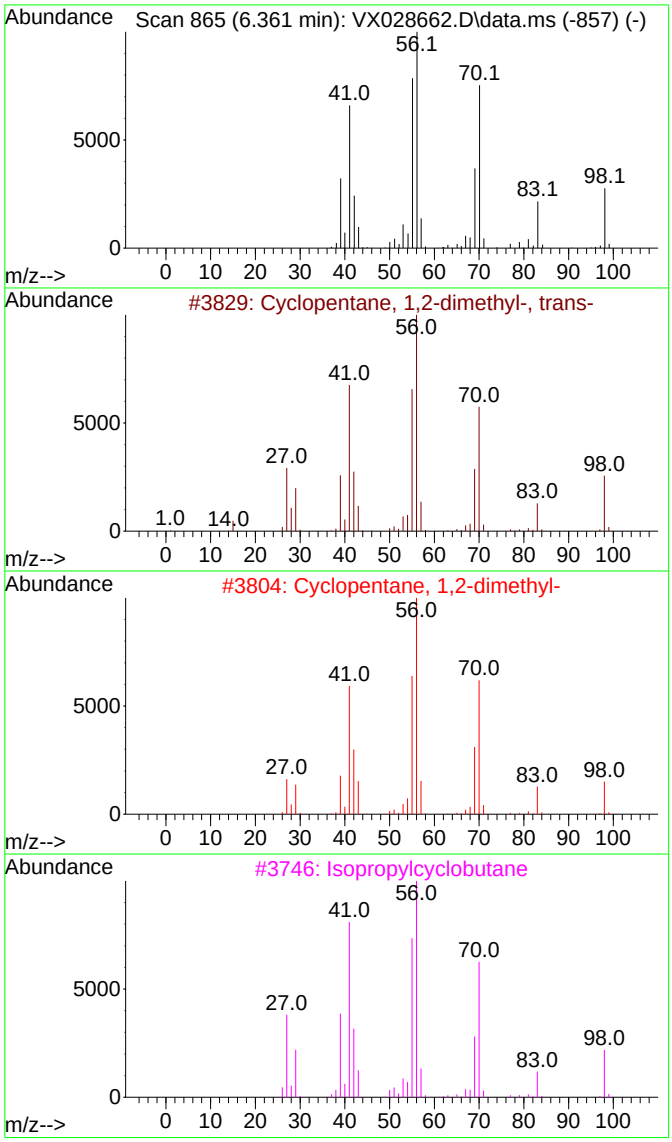
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

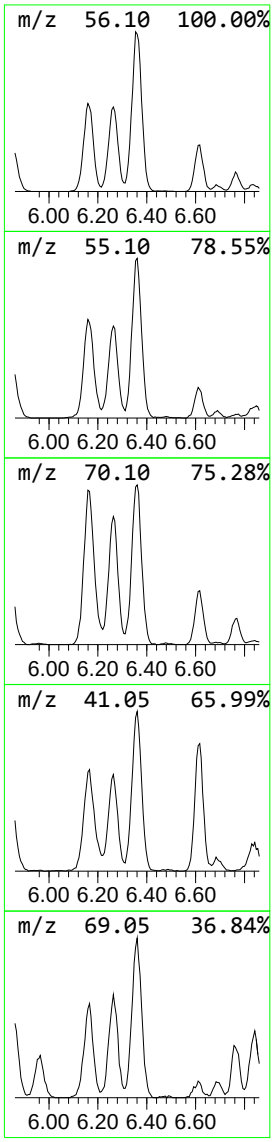
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Cyclopentane, 1,2-dimethyl-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.361	32.34 ug/l	659612	1,4-Difluorobenzene	6.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,2-dimethyl-, trans-	98	C7H14	000822-50-4	91
2	Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	91
3	Isopropylcyclobutane	98	C7H14	000872-56-0	90
4	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	86
5	Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	86





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
Data File : VX028662.D
Acq On : 13 May 2022 00:39
Operator : JC/MD
Sample : N2811-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

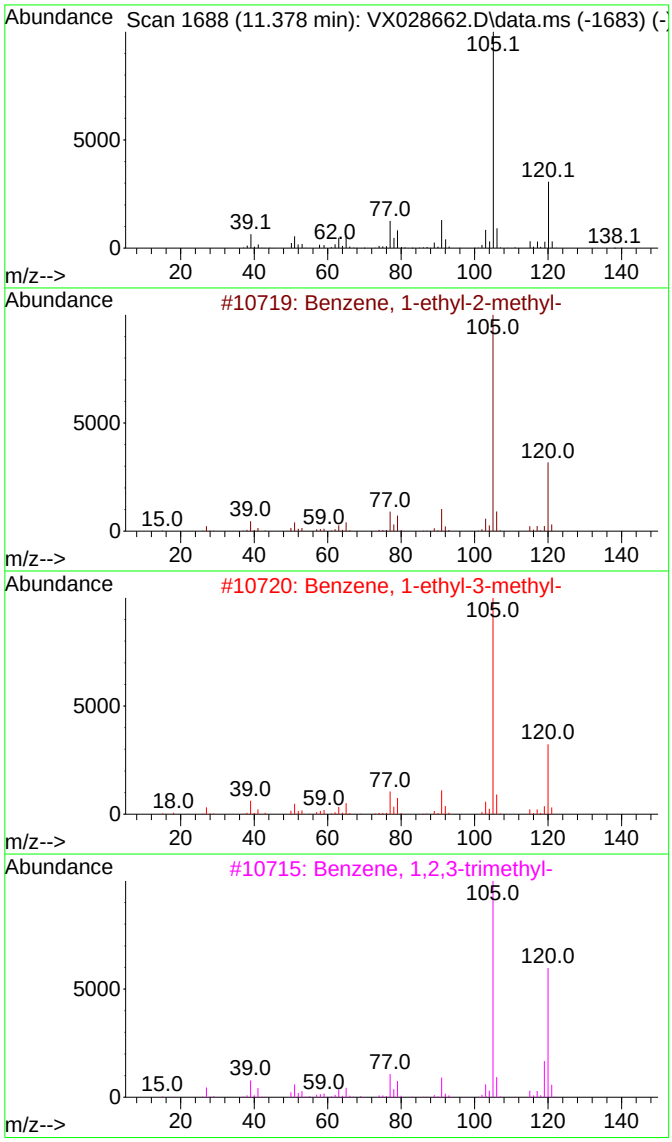
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

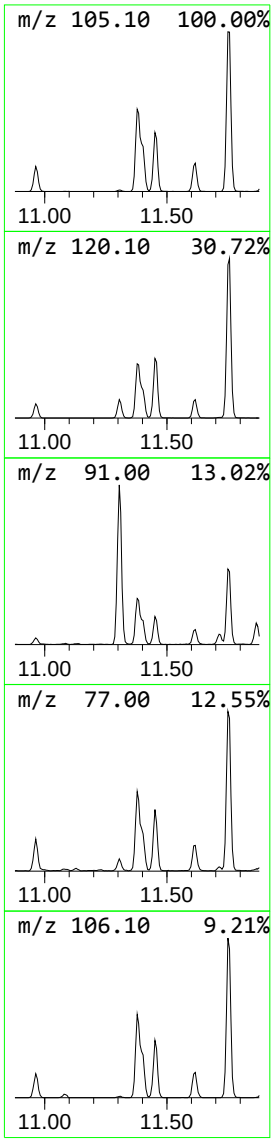
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 1-ethyl-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.378	46.04 ug/l	1289180	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-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
Data File : VX028662.D
Acq On : 13 May 2022 00:39
Operator : JC/MD
Sample : N2811-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.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	52.1	ug/l	501295	1	5.556	480754	50.0
Pentane	1.953	78.0	ug/l	749904	1	5.556	480754	50.0
Pentane, 2-methyl-	2.825	52.6	ug/l	505819	1	5.556	480754	50.0
Pentane, 3-methyl-	3.105	52.1	ug/l	500965	1	5.556	480754	50.0
n-Hexane	3.446	83.1	ug/l	798659	1	5.556	480754	50.0
Cyclopentane, m...	4.306	141.5	ug/l	1360870	1	5.556	480754	50.0
Hexane, 3-methyl-	5.836	54.4	ug/l	522854	1	5.556	480754	50.0
Cyclopentane, 1...	6.159	57.1	ug/l	549499	1	5.556	480754	50.0
Cyclopentane, 1...	6.361	32.3	ug/l	659612	2	6.763	1019700	50.0
Benzene, 1-ethy...	11.378	46.0	ug/l	1289180	4	12.024	1400060	50.0