LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062122\
Data File : VX029667.D
Acq On : 21 Jun 2022 17:19
Operator : JC/MD
Sample : N3290-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11ARE

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\82X061822W.M
Title : SW846 8260

Signal : TIC: VX029667.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.270	24	30	43	rBV5	7022	25742	1.89%	0.158%
2	1.752	101	109	115	rBV2	9402	14345	1.05%	0.088%
3	2.337	194	205	221	rVB	111177	226861	16.64%	1.391%
4	2.776	266	277	290	rBV	384806	822908	60.37%	5.047%
5	2.959	300	307	318	rVB2	16264	38771	2.84%	0.238%
6	3.111	325	332	342	rVB2	12230	29644	2.17%	0.182%
7	4.050	475	486	495	rBV3	7428	22388	1.64%	0.137%
8	4.209	497	512	525	rVV3	89838	273476	20.06%	1.677%
9	4.452	534	552	563	rBV5	12758	54101	3.97%	0.332%
10	5.123	649	662	679	rBV6	16699	70035	5.14%	0.430%
11	5.385	694	705	721	rBV	142593	421764	30.94%	2.587%
12	5.550	721	732	738	rVV	266382	776059	56.94%	4.759%
13	5.617	738	743	754	rVV	207551	640581	47.00%	3.929%
14	5.824	754	777	792	rVV4	114228	989114	72.57%	6.066%
15	5.952	792	798	816	rVB	158600	450743	33.07%	2.764%
16	6.245	828	846	857	rBV	171753	562101	41.24%	3.447%
17	6.354	857	864	875	rVB5	17858	51842	3.80%	0.318%
18	6.763	919	931	943	rBV	391419	962985	70.65%	5.906%
19	7.178	980	999	1021	rBV6	21926	84601	6.21%	0.519%
20	7.366	1021	1030	1042	rBB5	30456	81815	6.00%	0.502%
21	7.501	1045	1052	1056	rBV	17997	38199	2.80%	0.234%
22	7.562	1056	1062	1065	rVV3	27563	65655	4.82%	0.403%
23	7.592	1065	1067	1078	rVB	25456	55589	4.08%	0.341%
24	7.775	1090	1097	1103	rVB2	28232	59123	4.34%	0.363%
25	7.982	1117	1131	1140	rVB	146673	347037	25.46%	2.128%
26	8.110	1143	1152	1162	rBV	275280	576836	42.32%	3.538%
27	8.318	1179	1186	1194	rBV5	15033	41568	3.05%	0.255%
28	8.385	1194	1197	1200	rVV2	9785	15233	1.12%	0.093%
29	8.427	1200	1204	1210	rVB2	21843	37287	2.74%	0.229%
30	8.507	1210	1217	1226	rBV4	21023	47265	3.47%	0.290%
31	8.647	1228	1240	1252	rBV	805727	1363013	100.00%	8.359%
32	8.842	1264	1272	1276	rBV6	10453	24253	1.78%	0.149%
33	8.921	1282	1285	1289	rVV2	13692	19024	1.40%	0.117%
34	8.976	1289	1294	1299	rVV2	22638	35260	2.59%	0.216%
35	9.098	1307	1314	1319	rVB5	23497	48530	3.56%	0.298%
36	9.165	1319	1325	1334	rBV	93687	161134	11.82%	0.988%

LSC Area Percent Report

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Acq On : 21 Jun 2022 17:19
Operator : JC/MD
Sample : N3290-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11ARE

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\82X061822W.M
Title : SW846 8260

37	9.244	1334	1338	1344	rVB5	8261	14248	1.05%	0.087%
38	9.317	1344	1350	1356	rVB4	9961	19316	1.42%	0.118%
39	9.445	1364	1371	1377	rBV3	27138	64332	4.72%	0.395%
40	9.506	1377	1381	1387	rVB4	16336	30981	2.27%	0.190%
41	9.659	1402	1406	1412	rBV	12175	18259	1.34%	0.112%
42	9.762	1412	1423	1426	rBV6	27474	62542	4.59%	0.384%
43	10.000	1456	1462	1467	rBV	423239	773192	56.73%	4.742%
44	10.055	1467	1471	1476	rVV	808904	1063343	78.01%	6.521%
45	10.104	1476	1479	1482	rVV2	34685	49868	3.66%	0.306%
46	10.140	1482	1485	1491	rVB5	31945	58727	4.31%	0.360%
47	10.195	1491	1494	1498	rBV2	10454	19145	1.40%	0.117%
48	10.421	1526	1531	1536	rBV6	7982	14028	1.03%	0.086%
49	10.494	1536	1543	1548	rVB3	12850	21850	1.60%	0.134%
50	10.585	1554	1558	1561	rVV5	12059	18890	1.39%	0.116%
51	10.646	1563	1568	1573	rVB4	22034	38725	2.84%	0.237%
52	10.793	1587	1592	1598	rBV5	25800	62029	4.55%	0.380%
53	10.848	1598	1601	1610	rVB3	11667	24574	1.80%	0.151%
54	10.970	1614	1621	1625	rBV4	18560	33345	2.45%	0.205%
55	11.018	1625	1629	1631	rBV2	14412	18570	1.36%	0.114%
56	11.085	1635	1640	1644	rVB	606285	779696	57.20%	4.782%
57	11.177	1652	1655	1662	rVB3	15179	19307	1.42%	0.118%
58	11.268	1665	1670	1674	rVV3	16035	31397	2.30%	0.193%
59	11.311	1674	1677	1680	rVV3	13016	16653	1.22%	0.102%
60	11.372	1684	1687	1689	rVV	11392	15273	1.12%	0.094%
61	11.402	1689	1692	1695	rVV2	23183	34200	2.51%	0.210%
62	11.549	1707	1716	1725	rBV	660672	999076	73.30%	6.127%
63	11.701	1735	1741	1749	rBV7	15218	34731	2.55%	0.213%
64	11.817	1751	1760	1764	rBV4	21967	56320	4.13%	0.345%
65	11.890	1769	1772	1774	rVV	38870	55169	4.05%	0.338%
66	11.921	1774	1777	1783	rVB	258849	341598	25.06%	2.095%
67	12.024	1789	1794	1802	rVB	828234	1001902	73.51%	6.145%
68	12.110	1802	1808	1813	rBV6	21452	42628	3.13%	0.261%
69	12.177	1816	1819	1823	rVB	19970	27413	2.01%	0.168%
70	12.250	1823	1831	1838	rBV	292526	358862	26.33%	2.201%
71	12.445	1859	1863	1873	rVV7	9732	21999	1.61%	0.135%
72	12.524	1873	1876	1881	rVB3	27172	37572	2.76%	0.230%
73	12.628	1886	1893	1897	rBV	26441	37701	2.77%	0.231%
74	12.719	1902	1908	1912	rBV4	25623	42967	3.15%	0.264%
75	12.762	1912	1915	1920	rVB	15697	19830	1.45%	0.122%
76	12.841	1920	1928	1930	rBV	30182	47990	3.52%	0.294%

LSC Area Percent Report

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Acq On : 21 Jun 2022 17:19
Operator : JC/MD
Sample : N3290-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11ARE

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\82X061822W.M
Title : SW846 8260

77	12.866	1930	1932	1937	rVB3	26542	37533	2.75%	0.230%
78	13.036	1953	1960	1966	rVB4	52934	104348	7.66%	0.640%
79	13.195	1982	1986	1998	rVB	34382	54294	3.98%	0.333%
80	13.426	2016	2024	2030	rBV4	15528	33992	2.49%	0.208%
81	13.518	2035	2039	2047	rVB4	22523	46482	3.41%	0.285%
82	13.621	2051	2056	2061	rVV2	19389	31288	2.30%	0.192%
83	14.274	2156	2163	2167	rVV5	8721	13859	1.02%	0.085%
84	14.323	2167	2171	2177	rVV2	13406	20617	1.51%	0.126%

Sum of corrected areas: 16305543

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062122\
Data File : VX029667.D
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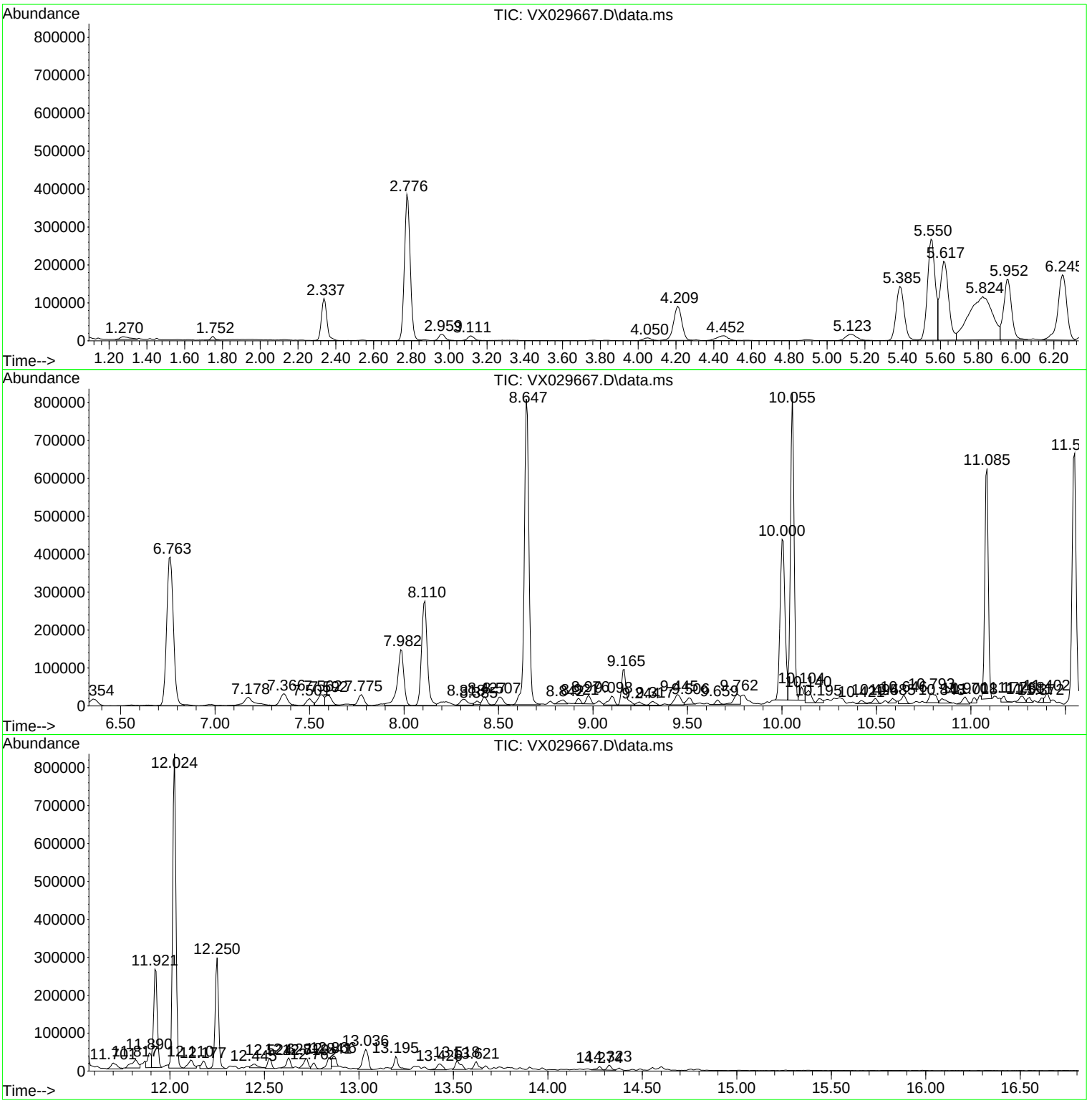
Operator : JC/MD
Sample : N3290-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11ARE

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062122\
Data File : VX029667.D
Acq On : 21 Jun 2022 17:19
Operator : JC/MD
Sample : N3290-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11ARE

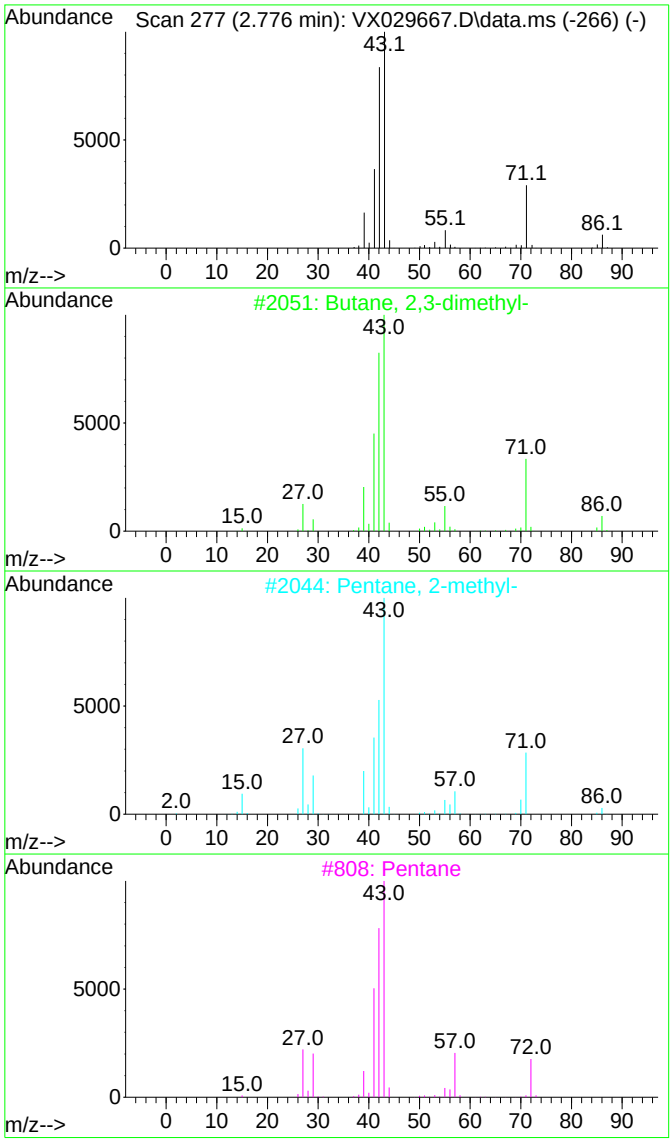
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

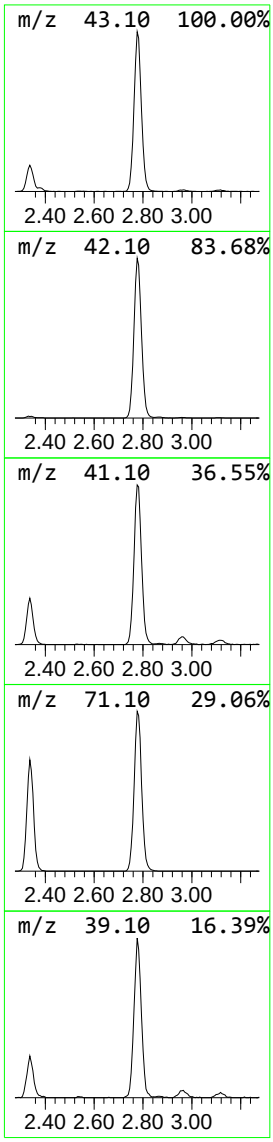
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Butane, 2,3-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.776	53.02 ug/l	822908	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	Pentane	72	C5H12	000109-66-0	38
4	Butanal, 2,2-dimethyl-	100	C6H12O	002094-75-9	12
5	C2H5C(CH3)2COCH3	114	C7H14O	020669-04-9	12





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062122\
Data File : VX029667.D
Acq On : 21 Jun 2022 17:19
Operator : JC/MD
Sample : N3290-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11ARE

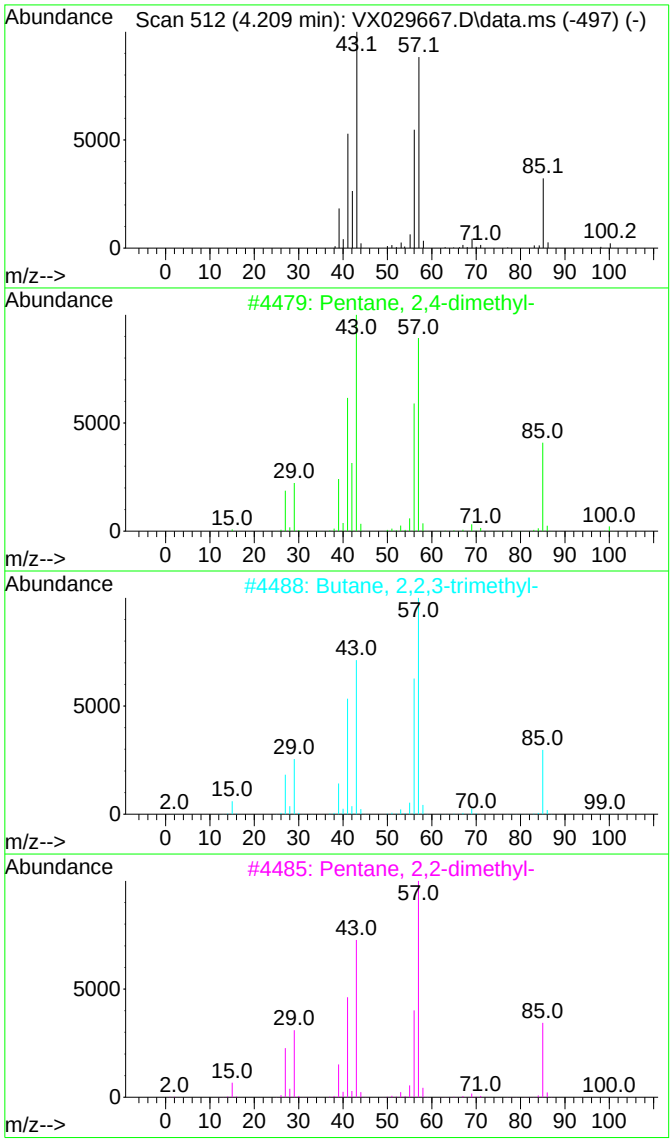
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

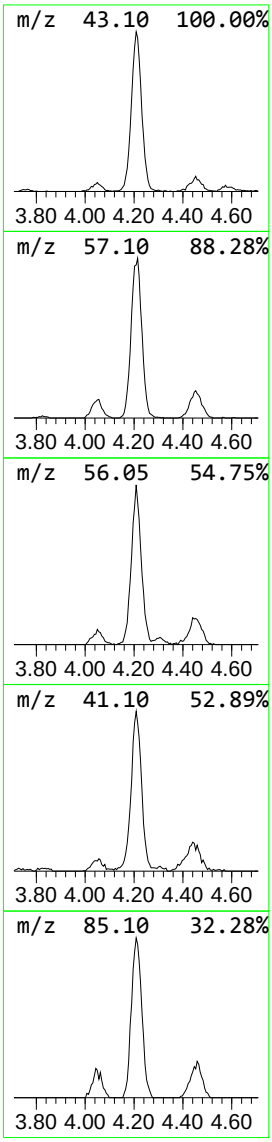
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Pentane, 2,4-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.209	17.62 ug/l	273476	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	94
2	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	59
3	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	47
4	Hexane, 2-methyl-	100	C7H16	000591-76-4	46
5	Pentane, 3-methyl-	86	C6H14	000096-14-0	43





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062122\
Data File : VX029667.D
Acq On : 21 Jun 2022 17:19
Operator : JC/MD
Sample : N3290-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11ARE

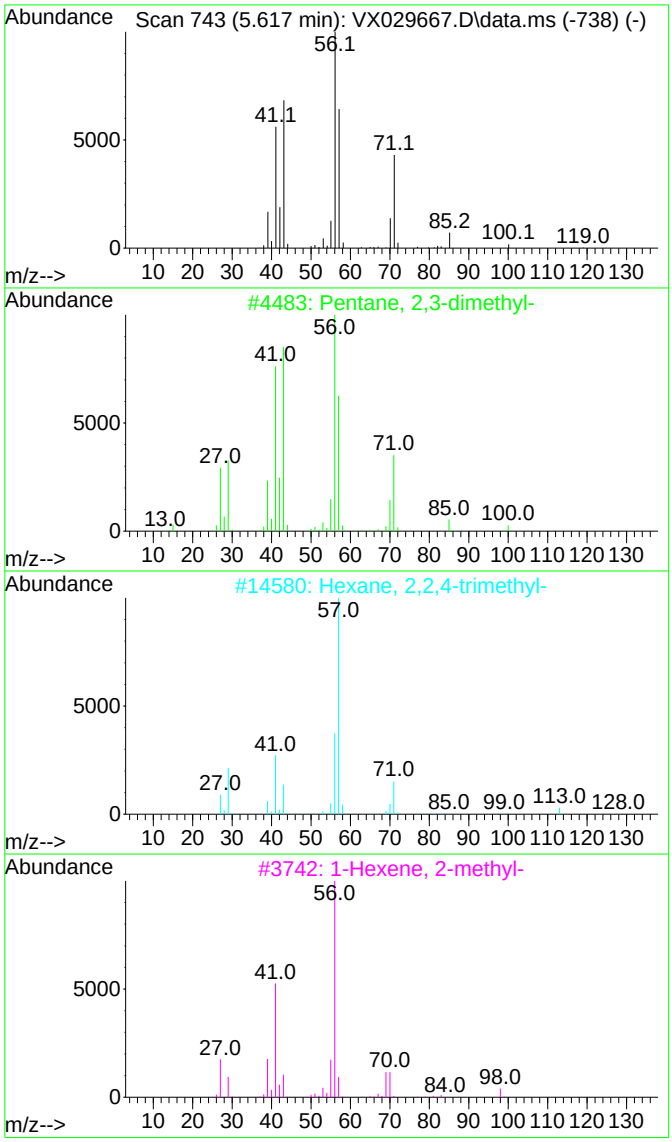
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

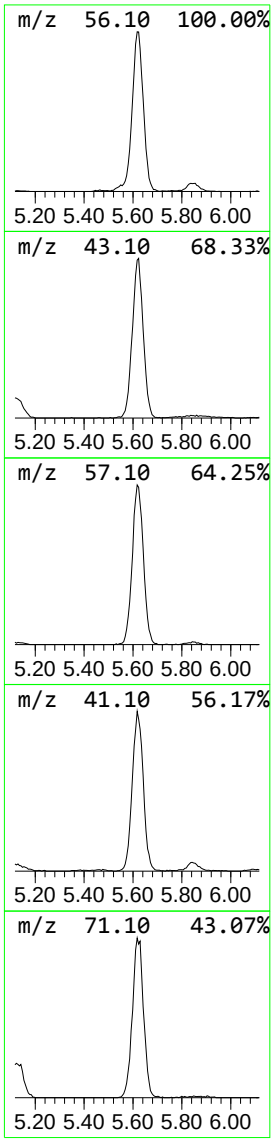
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Pentane, 2,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.617	41.27 ug/l	640581	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
2	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	39
3	1-Hexene, 2-methyl-	98	C7H14	006094-02-6	38
4	Heptane	100	C7H16	000142-82-5	37
5	Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	35





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062122\
Data File : VX029667.D
Acq On : 21 Jun 2022 17:19
Operator : JC/MD
Sample : N3290-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11ARE

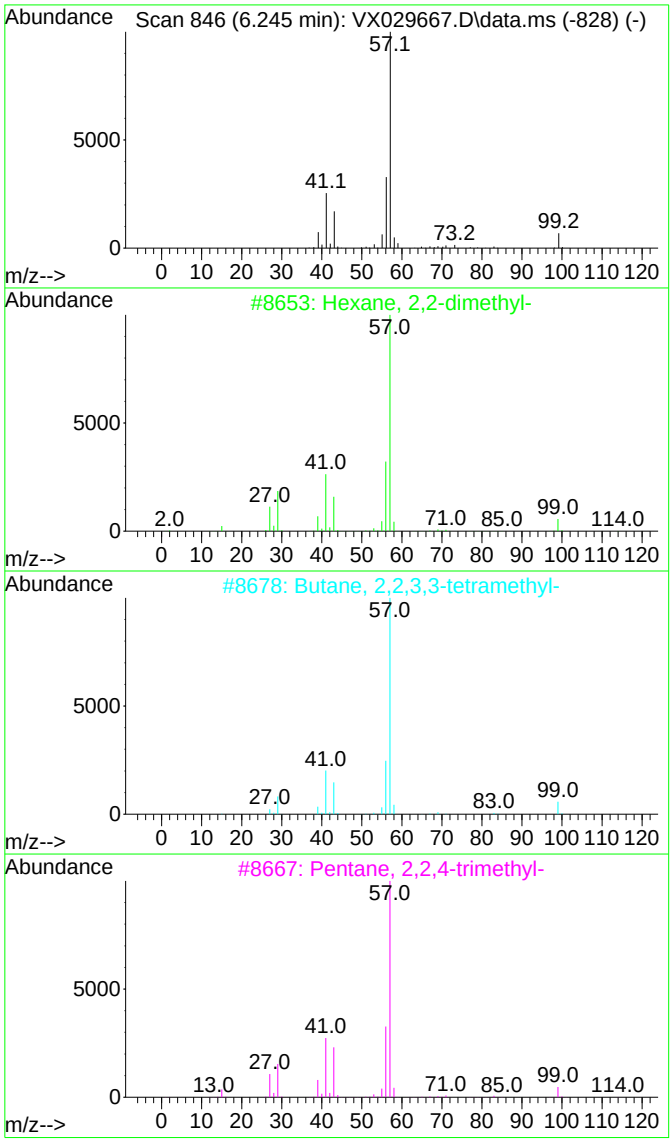
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

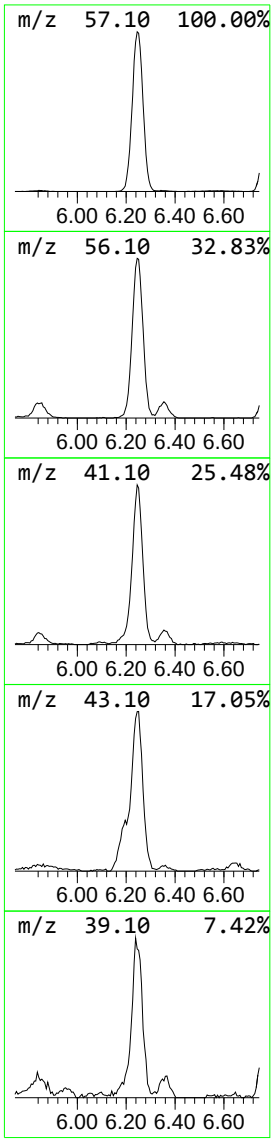
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Hexane, 2,2-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.245	29.19 ug/l	562101	1,4-Difluorobenzene	6.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	72
2	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	72
3	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
4	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	59
5	Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	50





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Data File : VX029667.D
Acq On : 21 Jun 2022 17:19
Operator : JC/MD
Sample : N3290-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11ARE

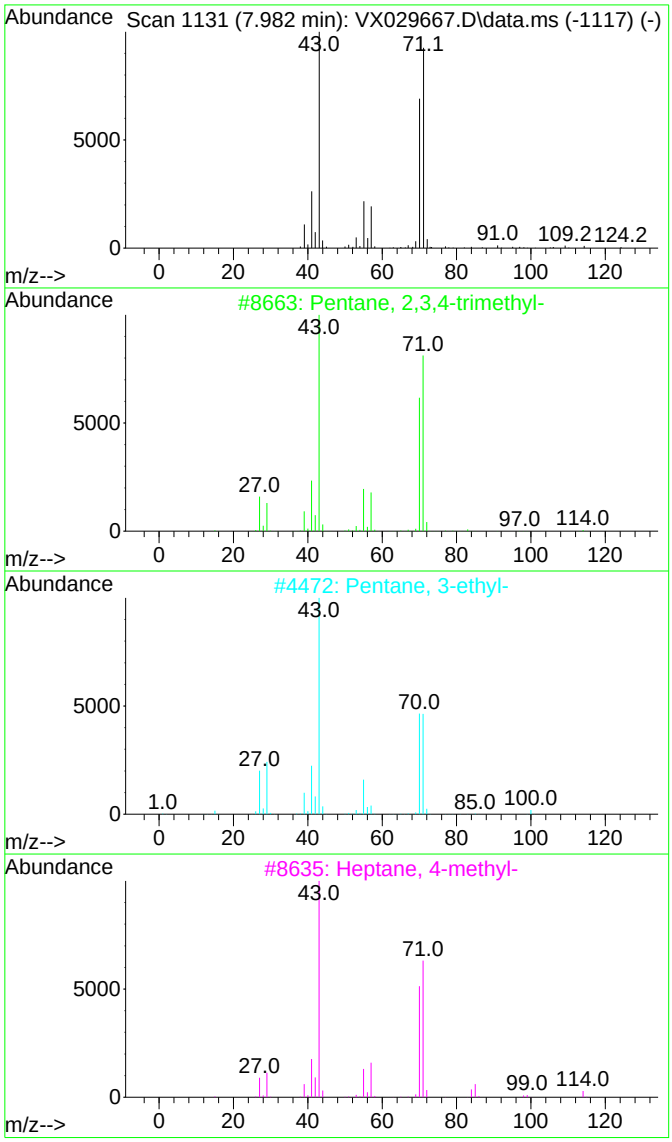
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

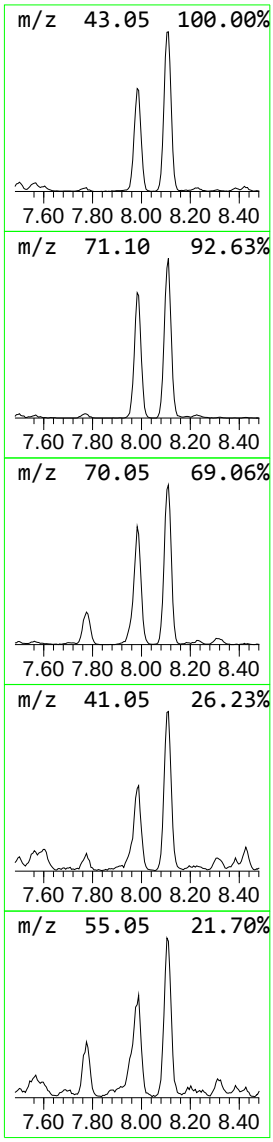
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Pentane, 2,3,4-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.982	18.02 ug/l	347037	1,4-Difluorobenzene	6.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	91
2	Pentane, 3-ethyl-	100	C7H16	000617-78-7	83
3	Heptane, 4-methyl-	114	C8H18	000589-53-7	74
4	Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	64
5	Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062122\
Data File : VX029667.D
Acq On : 21 Jun 2022 17:19
Operator : JC/MD
Sample : N3290-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11ARE

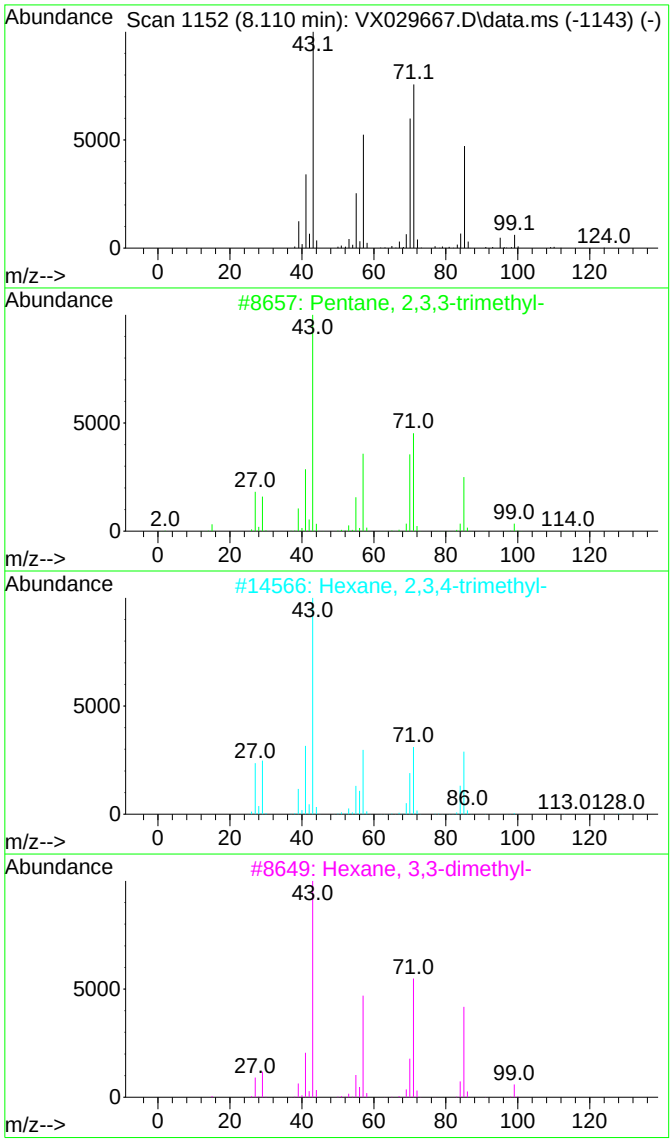
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

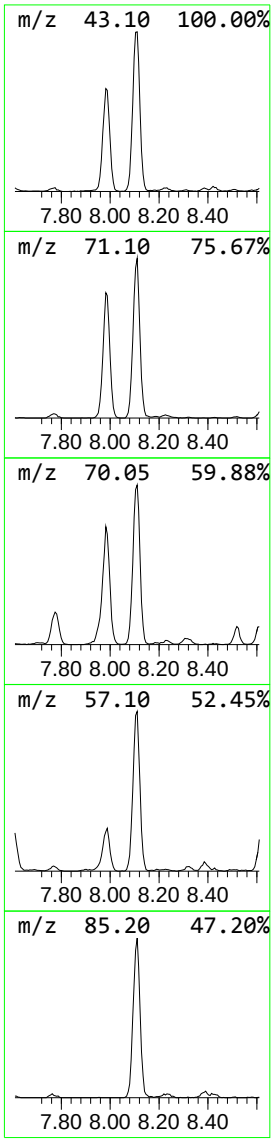
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Pentane, 2,3,3-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.110	29.95 ug/l	576836	1,4-Difluorobenzene	6.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	64
3	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	64
4	Heptane, 4-methyl-	114	C8H18	000589-53-7	59
5	Hexane, 3-methyl-	100	C7H16	000589-34-4	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062122\
Data File : VX029667.D
Acq On : 21 Jun 2022 17:19
Operator : JC/MD
Sample : N3290-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11ARE

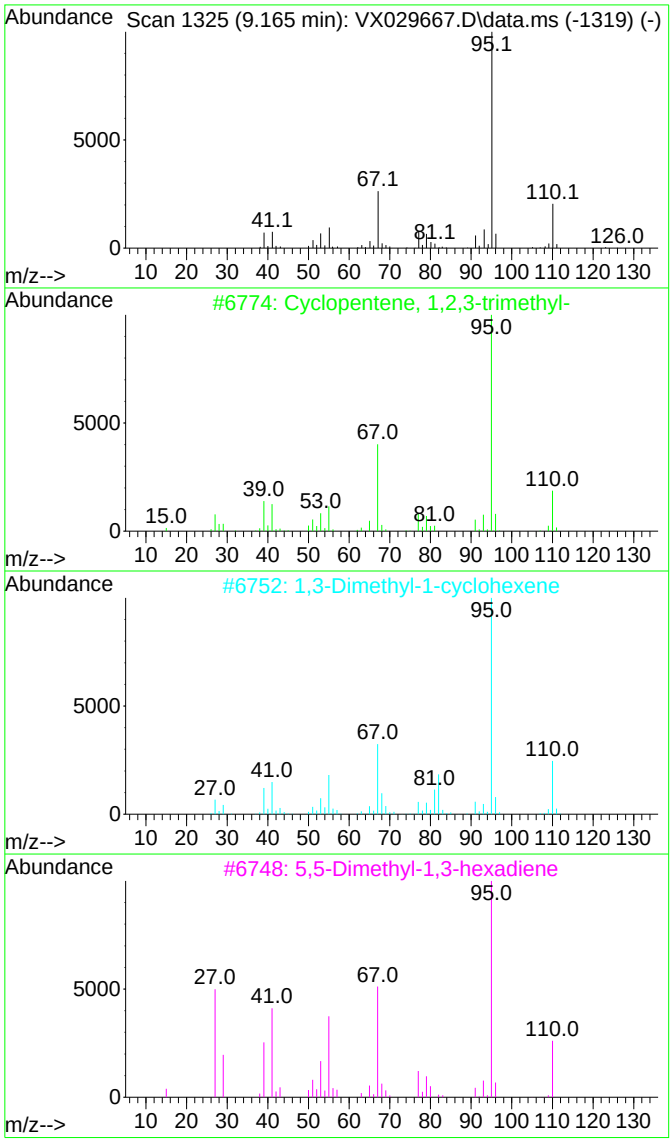
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

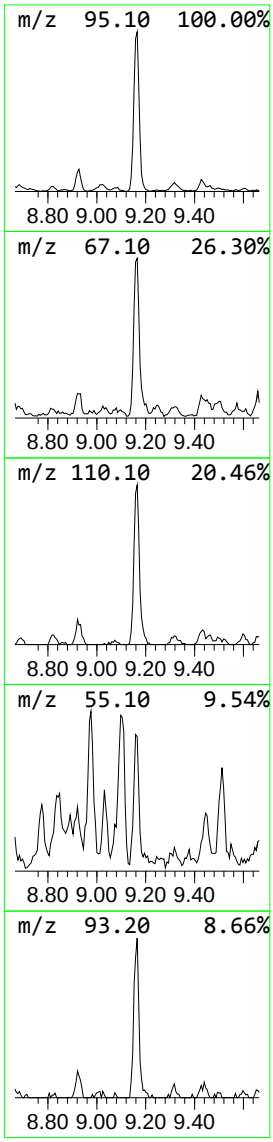
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Cyclopentene, 1,2,3-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.165	7.58 ug/l	161134	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentene, 1,2,3-trimethyl-	110	C8H14	000473-91-6	91
2	1,3-Dimethyl-1-cyclohexene	110	C8H14	002808-76-6	91
3	5,5-Dimethyl-1,3-hexadiene	110	C8H14	001515-79-3	87
4	Bicyclo[3.1.0]hexane, 1,5-dimethyl-	110	C8H14	1010142-17-5	72
5	Bicyclo[2.2.1]heptane, 2-(2-prop...	136	C10H16	002633-80-9	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062122\
Data File : VX029667.D
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Operator : JC/MD
Sample : N3290-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11ARE

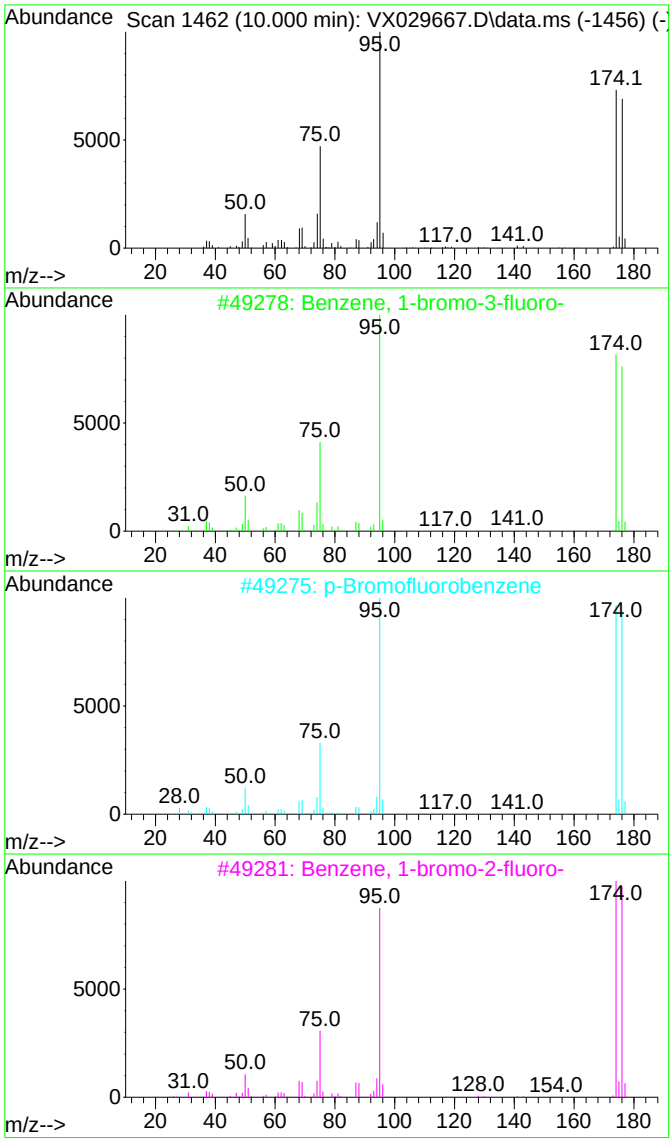
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

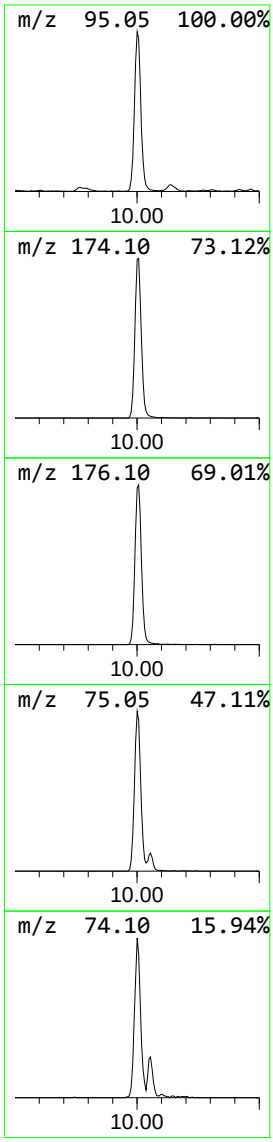
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1-bromo-3-fluoro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.000	36.36 ug/l	773192	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-bromo-3-fluoro-	174	C6H4BrF	001073-06-9	94
2	p-Bromofluorobenzene	174	C6H4BrF	000460-00-4	93
3	Benzene, 1-bromo-2-fluoro-	174	C6H4BrF	001072-85-1	91
4	2(1H)-Pyrimidinone, 5-bromo-	174	C4H3BrN2O	038353-06-9	36
5	4(1H)-Pyrimidinone, 5-bromo-	174	C4H3BrN2O	019808-30-1	32





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062122\
Data File : VX029667.D
Acq On : 21 Jun 2022 17:19
Operator : JC/MD
Sample : N3290-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11ARE

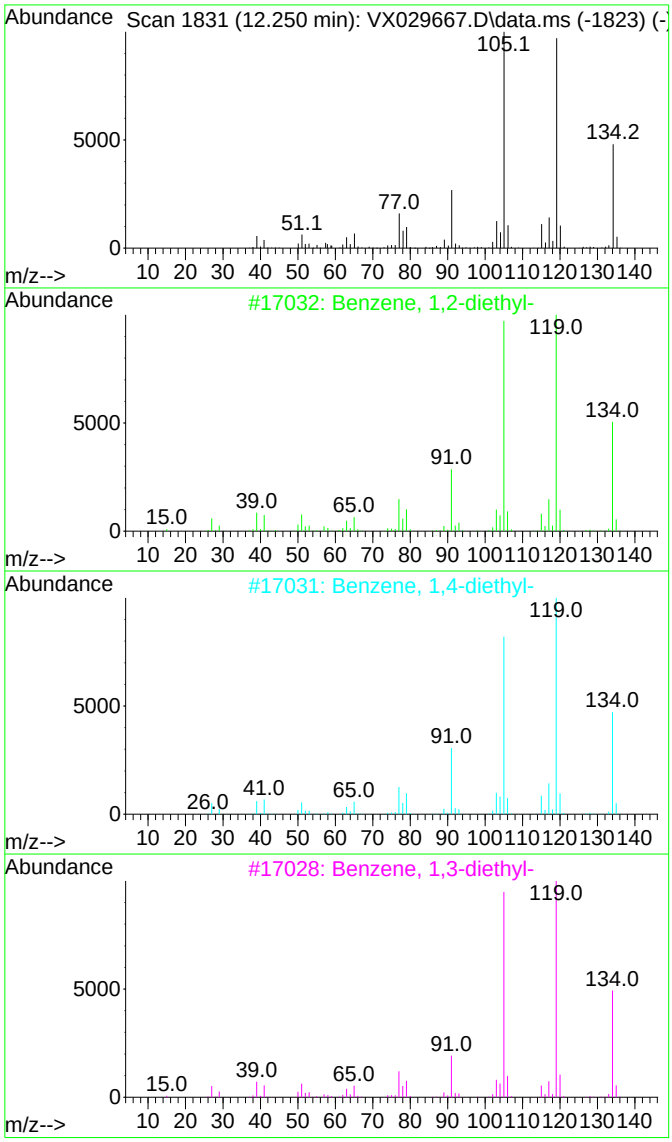
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

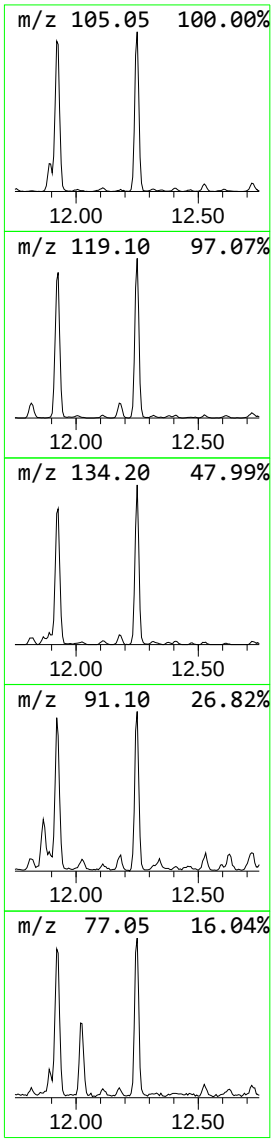
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Benzene, 1,2-diethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.250	17.91 ug/l	358862	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	97
2	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	96
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	94
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91





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Data File : VX029667.D
Acq On : 21 Jun 2022 17:19
Operator : JC/MD
Sample : N3290-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11ARE

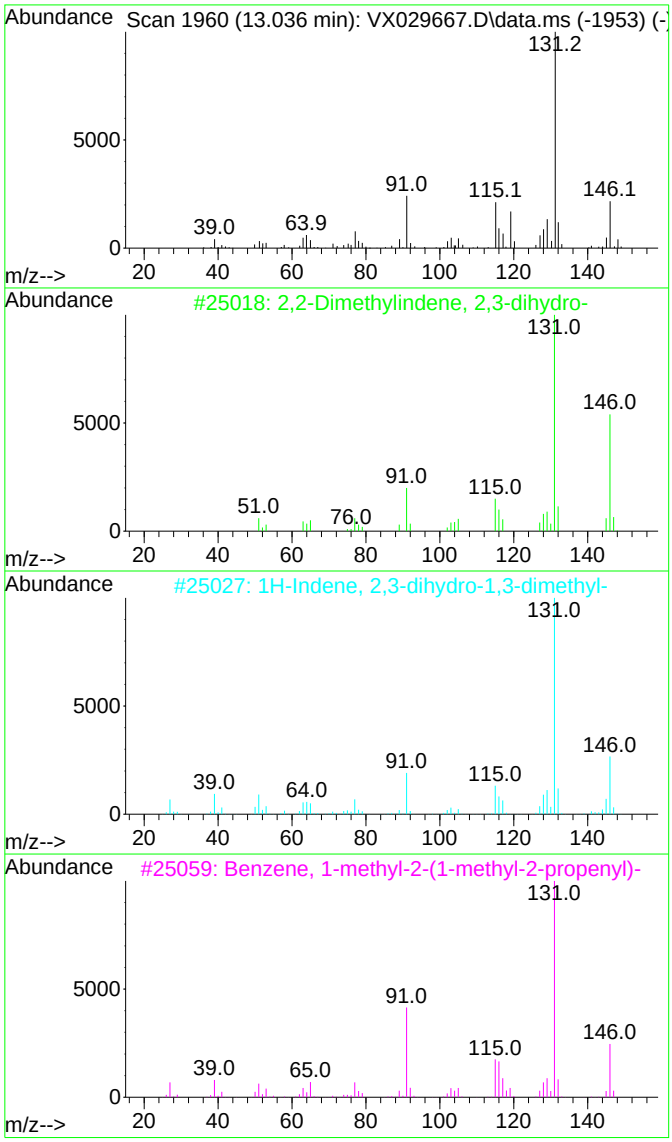
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

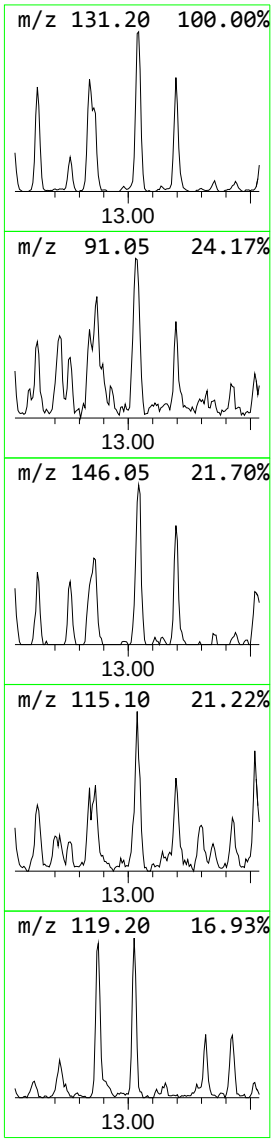
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 2,2-Dimethylindene, 2,3-dih... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.036	5.21 ug/l	104348	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	87
2	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	87
3	Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	83
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	83
5	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	83





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062122\
Data File : VX029667.D
Acq On : 21 Jun 2022 17:19
Operator : JC/MD
Sample : N3290-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11ARE

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.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, 2,3-dim...	2.776	53.0	ug/l	822908	1	5.556	776059	50.0
Pentane, 2,4-di...	4.209	17.6	ug/l	273476	1	5.556	776059	50.0
Pentane, 2,3-di...	5.617	41.3	ug/l	640581	1	5.556	776059	50.0
Hexane, 2,2-dim...	6.245	29.2	ug/l	562101	2	6.763	962985	50.0
Pentane, 2,3,4-...	7.982	18.0	ug/l	347037	2	6.763	962985	50.0
Pentane, 2,3,3-...	8.110	29.9	ug/l	576836	2	6.763	962985	50.0
Cyclopentene, 1...	9.165	7.6	ug/l	161134	3	10.055	1063340	50.0
Benzene, 1-brom...	10.000	36.4	ug/l	773192	3	10.055	1063340	50.0
Benzene, 1,2-di...	12.250	17.9	ug/l	358862	4	12.024	1001900	50.0
2,2-Dimethylind...	13.036	5.2	ug/l	104348	4	12.024	1001900	50.0