LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028103.D
Acq On : 14 Apr 2022 13:41
Operator : JC/MD
Sample : N2403-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-29R

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: VX028103.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.563	70	78	85	rBV4	17656	44196	1.21%	0.141%
2	1.752	104	109	116	rBV	229301	316239	8.69%	1.006%
3	2.337	198	205	218	rVB	36001	77544	2.13%	0.247%
4	2.782	270	278	289	rBV	134068	329017	9.04%	1.046%
5	2.959	299	307	321	rVV	159146	352319	9.68%	1.121%
6	3.111	321	332	346	rVB	657757	1523889	41.87%	4.847%
7	3.763	426	439	449	rBV3	26485	82853	2.28%	0.264%
8	4.215	501	513	521	rBV3	23728	71499	1.96%	0.227%
9	4.306	521	528	539	rVB2	20062	56906	1.56%	0.181%
10	5.391	693	706	715	rBV	150528	435862	11.98%	1.386%
11	5.556	724	733	741	rBV	224058	646366	17.76%	2.056%
12	5.623	741	744	760	rVB2	68930	173368	4.76%	0.551%
13	5.842	768	780	790	rBV5	20073	67714	1.86%	0.215%
14	5.958	790	799	807	rVV	157263	418578	11.50%	1.331%
15	6.044	807	813	826	rVV2	64379	197249	5.42%	0.627%
16	6.165	826	833	835	rVV5	31009	75341	2.07%	0.240%
17	6.239	835	845	859	rVV	1294480	3639203	100.00%	11.575%
18	6.354	859	864	877	rVB5	16264	51118	1.40%	0.163%
19	6.763	921	931	941	rBV	380112	905205	24.87%	2.879%
20	7.373	1021	1031	1044	rVB5	29689	86749	2.38%	0.276%
21	7.988	1123	1132	1144	rVB2	36091	94845	2.61%	0.302%
22	8.110	1144	1152	1161	rBV	79810	174347	4.79%	0.555%
23	8.427	1195	1204	1210	rVV	451450	753862	20.72%	2.398%
24	8.507	1210	1217	1227	rVV	1103903	1777499	48.84%	5.653%
25	8.604	1227	1233	1235	rVV2	25322	42778	1.18%	0.136%
26	8.653	1235	1241	1251	rVB	827877	1317452	36.20%	4.190%
27	8.775	1251	1261	1264	rBV4	22967	54584	1.50%	0.174%
28	8.842	1264	1272	1280	rVV	1122186	1724706	47.39%	5.485%
29	8.976	1288	1294	1301	rVV4	26933	72568	1.99%	0.231%
30	9.104	1311	1315	1323	rVB2	23728	41487	1.14%	0.132%
31	9.445	1364	1371	1378	rBV2	218501	360062	9.89%	1.145%
32	9.561	1386	1390	1395	rVB3	48617	72548	1.99%	0.231%
33	10.025	1459	1466	1468	rVV	556153	838585	23.04%	2.667%
34	10.055	1468	1471	1476	rVV	894310	1169929	32.15%	3.721%
35	10.104	1476	1479	1482	rVV	213984	288412	7.93%	0.917%
36	10.146	1482	1486	1491	rVV	1006925	1309862	35.99%	4.166%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028103.D
Acq On : 14 Apr 2022 13:41
Operator : JC/MD
Sample : N2403-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-29R

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	10.195	1491	1494	1497	rVV	55454	70829	1.95%	0.225%
38	10.250	1497	1503	1509	rVV5	34320	85515	2.35%	0.272%
39	10.354	1516	1520	1528	rVV2	104019	174623	4.80%	0.555%
40	10.646	1563	1568	1574	rVB	62994	85446	2.35%	0.272%
41	10.793	1584	1592	1598	rBV2	232297	444309	12.21%	1.413%
42	10.848	1598	1601	1605	rVV	33074	51659	1.42%	0.164%
43	10.902	1605	1610	1614	rVV	91940	133571	3.67%	0.425%
44	10.963	1614	1620	1625	rVV3	51273	92039	2.53%	0.293%
45	11.085	1633	1640	1644	rVV	747232	1020447	28.04%	3.246%
46	11.122	1644	1646	1648	rVV	74059	85952	2.36%	0.273%
47	11.171	1648	1654	1661	rVV2	330106	662699	18.21%	2.108%
48	11.274	1665	1671	1674	rVV	523341	664779	18.27%	2.114%
49	11.311	1674	1677	1685	rVV2	195300	282858	7.77%	0.900%
50	11.390	1685	1690	1694	rVV2	113347	165562	4.55%	0.527%
51	11.439	1694	1698	1709	rVB3	78612	163716	4.50%	0.521%
52	11.616	1720	1727	1730	rBV	29128	45199	1.24%	0.144%
53	11.695	1734	1740	1746	rVV4	87243	163749	4.50%	0.521%
54	11.756	1746	1750	1753	rVV	22630	40941	1.12%	0.130%
55	11.799	1753	1757	1762	rVV	136315	178174	4.90%	0.567%
56	11.890	1762	1772	1778	rVV5	232448	455979	12.53%	1.450%
57	11.939	1778	1780	1789	rVV5	31517	70230	1.93%	0.223%
58	12.024	1789	1794	1800	rVV	949834	1191648	32.74%	3.790%
59	12.110	1800	1808	1812	rVV2	194187	371071	10.20%	1.180%
60	12.152	1812	1815	1817	rVV2	53069	81516	2.24%	0.259%
61	12.177	1817	1819	1824	rVV3	44346	55787	1.53%	0.177%
62	12.244	1824	1830	1838	rVV3	848522	1295859	35.61%	4.121%
63	12.317	1838	1842	1846	rVV	83422	120711	3.32%	0.384%
64	12.408	1850	1857	1862	rVV2	38178	73764	2.03%	0.235%
65	12.463	1862	1866	1872	rVV3	31342	65851	1.81%	0.209%
66	12.518	1872	1875	1879	rVV2	36218	51936	1.43%	0.165%
67	12.609	1885	1890	1893	rVV2	86787	157473	4.33%	0.501%
68	12.646	1893	1896	1901	rVV	248122	336641	9.25%	1.071%
69	12.701	1901	1905	1912	rVV	519071	707542	19.44%	2.250%
70	12.762	1912	1915	1922	rVV3	62881	117897	3.24%	0.375%
71	12.835	1922	1927	1935	rVV2	65208	119299	3.28%	0.379%
72	12.920	1935	1941	1944	rVV	72405	121076	3.33%	0.385%
73	12.963	1945	1948	1953	rVV2	63985	93581	2.57%	0.298%
74	13.030	1953	1959	1965	rVV5	38813	86123	2.37%	0.274%
75	13.097	1967	1970	1974	rVV3	31578	47653	1.31%	0.152%
76	13.134	1974	1976	1979	rVV	33240	43080	1.18%	0.137%

LSC Area Percent Report

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Acq On : 14 Apr 2022 13:41
Operator : JC/MD
Sample : N2403-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-29R

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	13.170	1979	1982	1984	rVV	38694	49134	1.35%	0.156%
78	13.292	1998	2002	2009	rVV	309287	438955	12.06%	1.396%
79	13.347	2009	2011	2017	rVV3	30937	60747	1.67%	0.193%
80	13.420	2020	2023	2030	rVV2	25440	46090	1.27%	0.147%
81	13.506	2033	2037	2040	rVV3	24064	39532	1.09%	0.126%
82	13.542	2040	2043	2047	rVV	58280	81666	2.24%	0.260%
83	13.585	2047	2050	2054	rVV2	68061	111307	3.06%	0.354%
84	13.640	2054	2059	2061	rVV2	76082	126218	3.47%	0.401%
85	13.664	2061	2063	2069	rVB2	91023	109089	3.00%	0.347%
86	13.780	2079	2082	2090	rVB	25382	39442	1.08%	0.125%
87	13.853	2090	2094	2101	rBV2	32320	45807	1.26%	0.146%
88	13.926	2101	2106	2110	rBV2	66725	88884	2.44%	0.283%
89	14.079	2126	2131	2134	rBV	35739	46797	1.29%	0.149%
90	14.213	2148	2153	2160	rBV	49881	77169	2.12%	0.245%
91	14.371	2175	2179	2184	rVB2	34937	47782	1.31%	0.152%
92	14.481	2192	2197	2203	rBV2	47934	66542	1.83%	0.212%
93	14.786	2242	2247	2251	rBV3	24673	44789	1.23%	0.142%

Sum of corrected areas: 31441475

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Data File : VX028103.D
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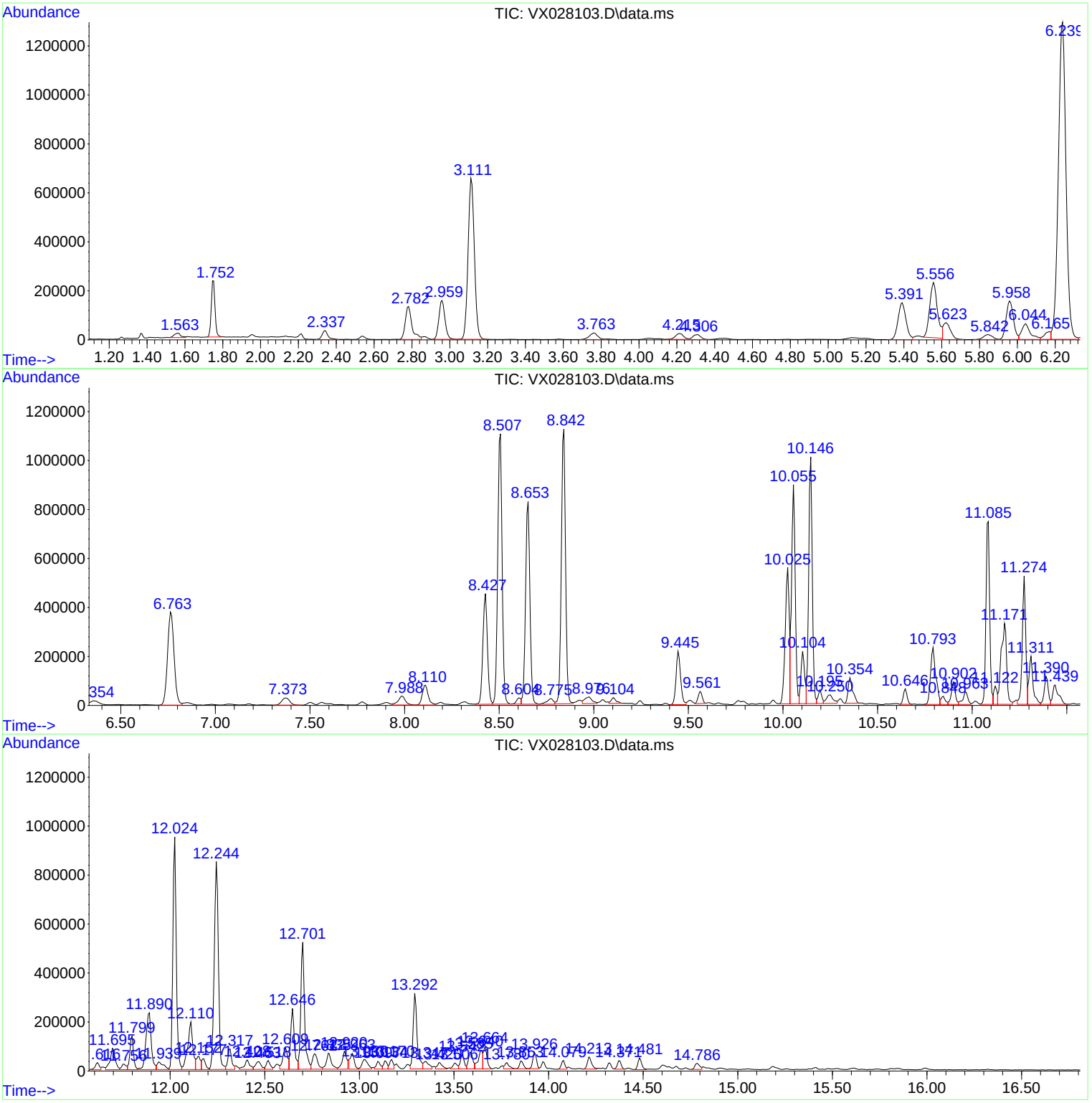
Sample : N2403-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-29R

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 : VX028103.D
Acq On : 14 Apr 2022 13:41
Operator : JC/MD
Sample : N2403-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-29R

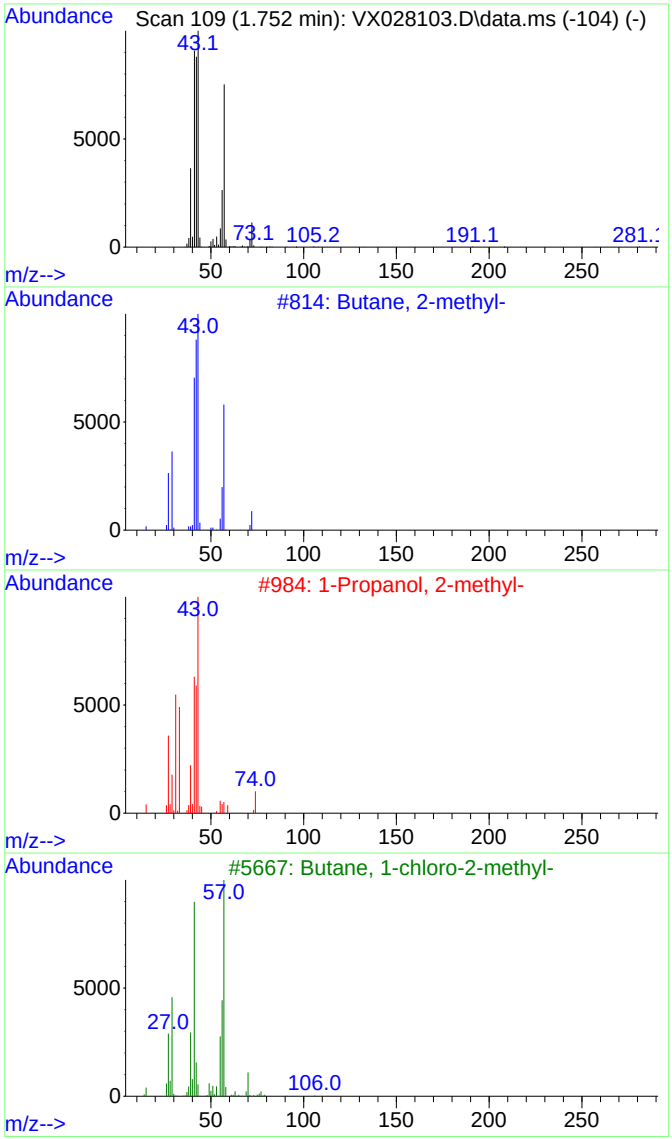
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

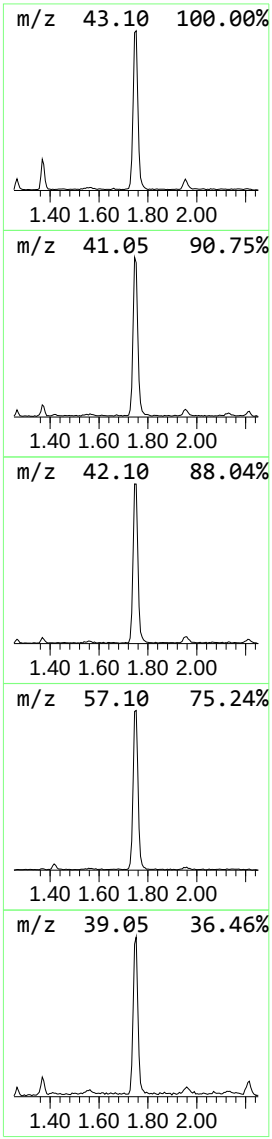
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Butane, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.752	24.46 ug/l	316239	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	86
2	1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	36
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	33





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028103.D
Acq On : 14 Apr 2022 13:41
Operator : JC/MD
Sample : N2403-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-29R

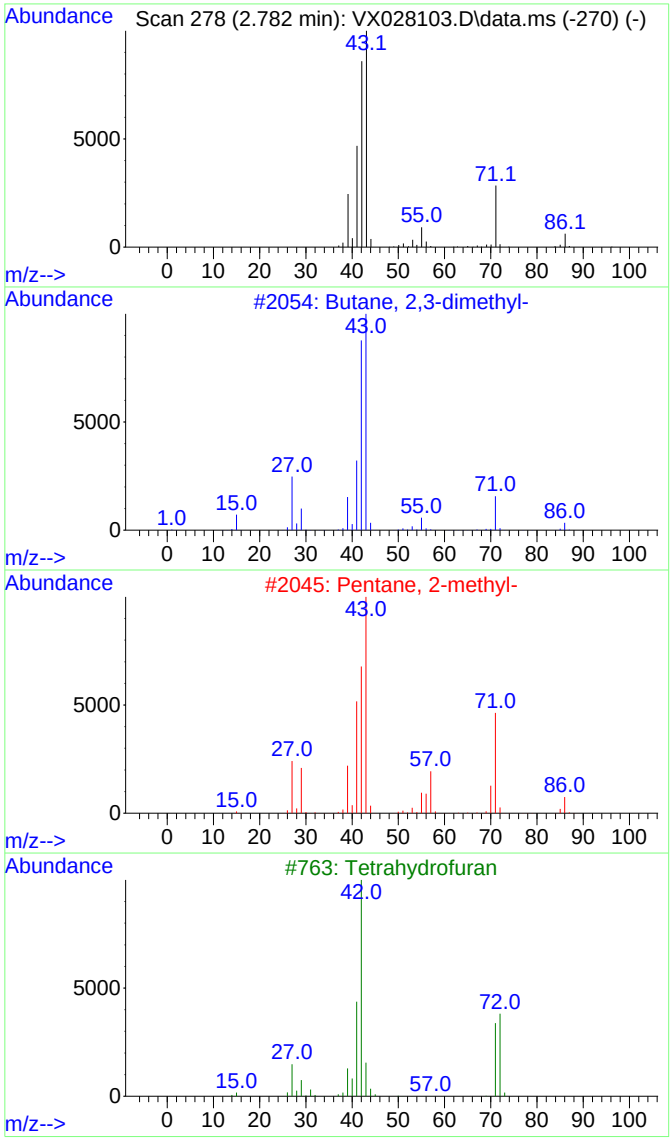
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

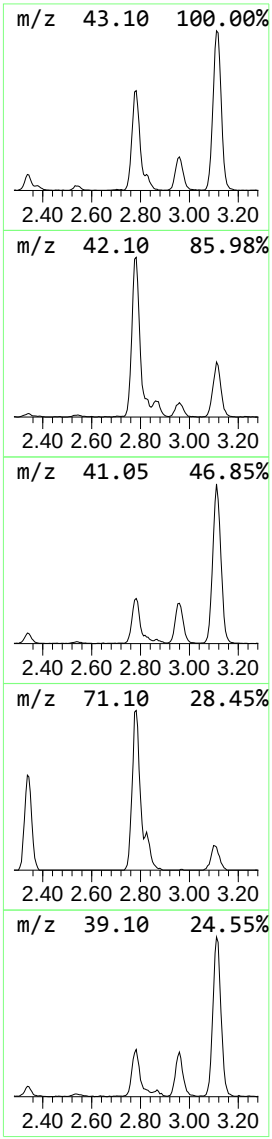
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Butane, 2,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.782	25.45 ug/l	329017	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	86
2	Pentane, 2-methyl-	86	C6H14	000107-83-5	78
3	Tetrahydrofuran	72	C4H8O	000109-99-9	38
4	Pentane	72	C5H12	000109-66-0	36
5	1-Pentene	70	C5H10	000109-67-1	28





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028103.D
Acq On : 14 Apr 2022 13:41
Operator : JC/MD
Sample : N2403-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-29R

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

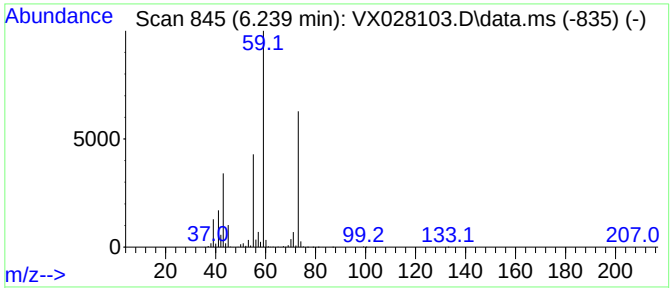
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

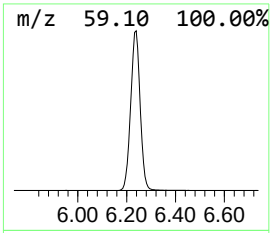
****

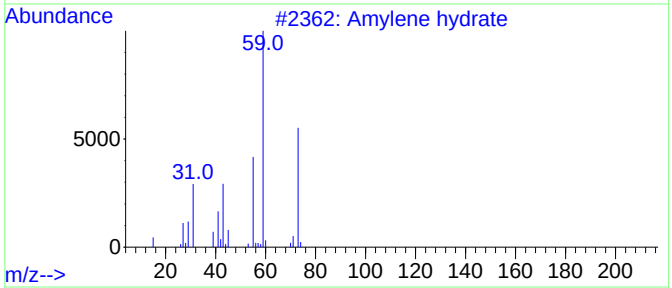
Peak Number 3 Amylene hydrate Concentration Rank 1

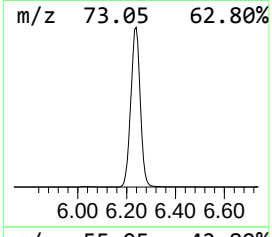
R.T.	EstConc	Area	Relative to ISTD	R.T.
6.239	201.01 ug/l	3639200	1,4-Difluorobenzene	6.763

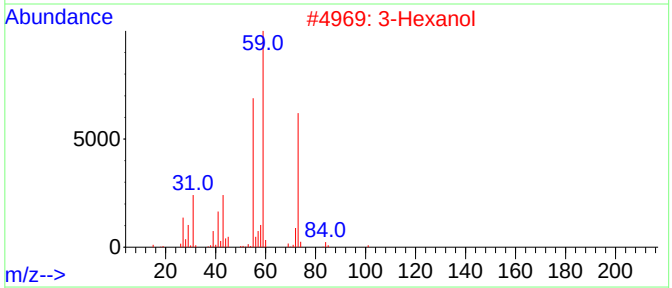
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Amylene hydrate	88	C5H12O	000075-85-4	83
2	3-Hexanol	102	C6H14O	000623-37-0	72
3	3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	40
4	3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	39
5	Silane, trimethyl-	74	C3H10Si	000993-07-7	38

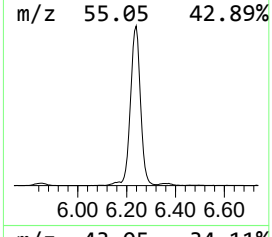


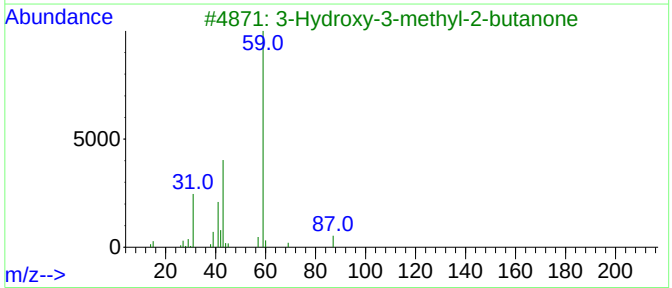


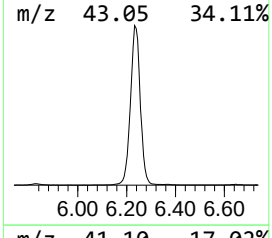


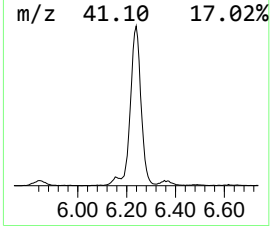












Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028103.D
Acq On : 14 Apr 2022 13:41
Operator : JC/MD
Sample : N2403-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-29R

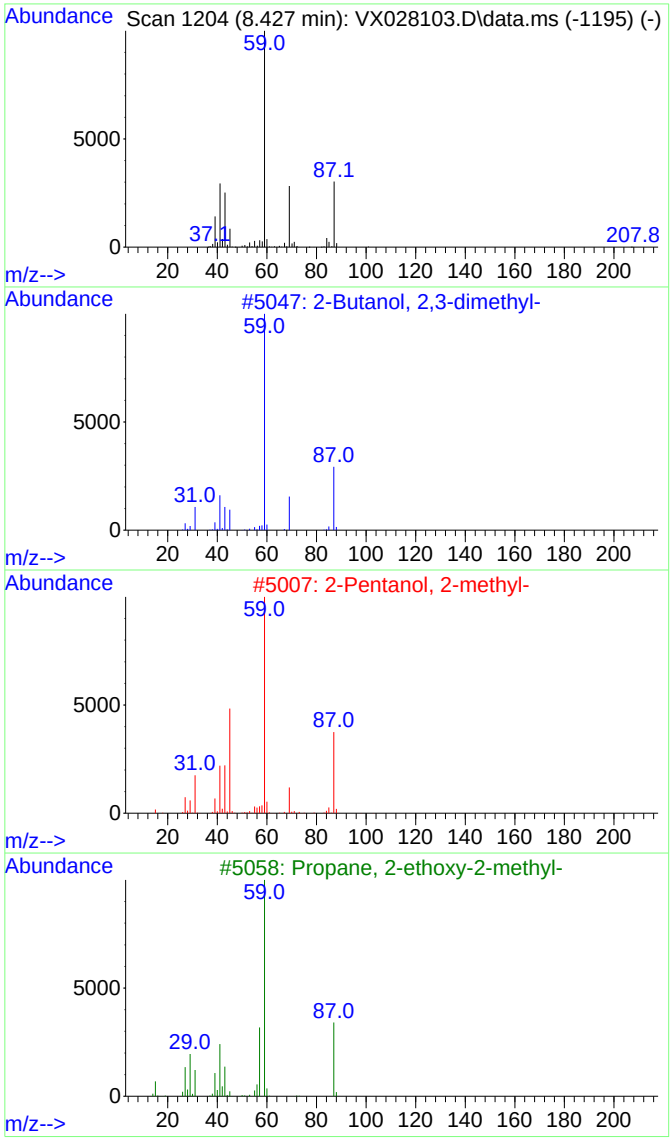
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

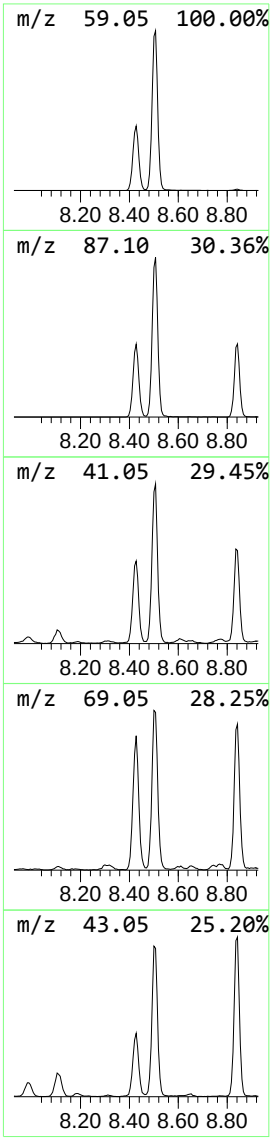
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 2-Butanol, 2,3-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.427	32.22 ug/l	753862	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanol, 2,3-dimethyl-	102	C6H14O	000594-60-5	83
2	2-Pentanol, 2-methyl-	102	C6H14O	000590-36-3	78
3	Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	56
4	3-Pentanol, 2,2-dimethyl-	116	C7H16O	003970-62-5	43
5	3-Heptanol	116	C7H16O	000589-82-2	40





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028103.D
Acq On : 14 Apr 2022 13:41
Operator : JC/MD
Sample : N2403-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-29R

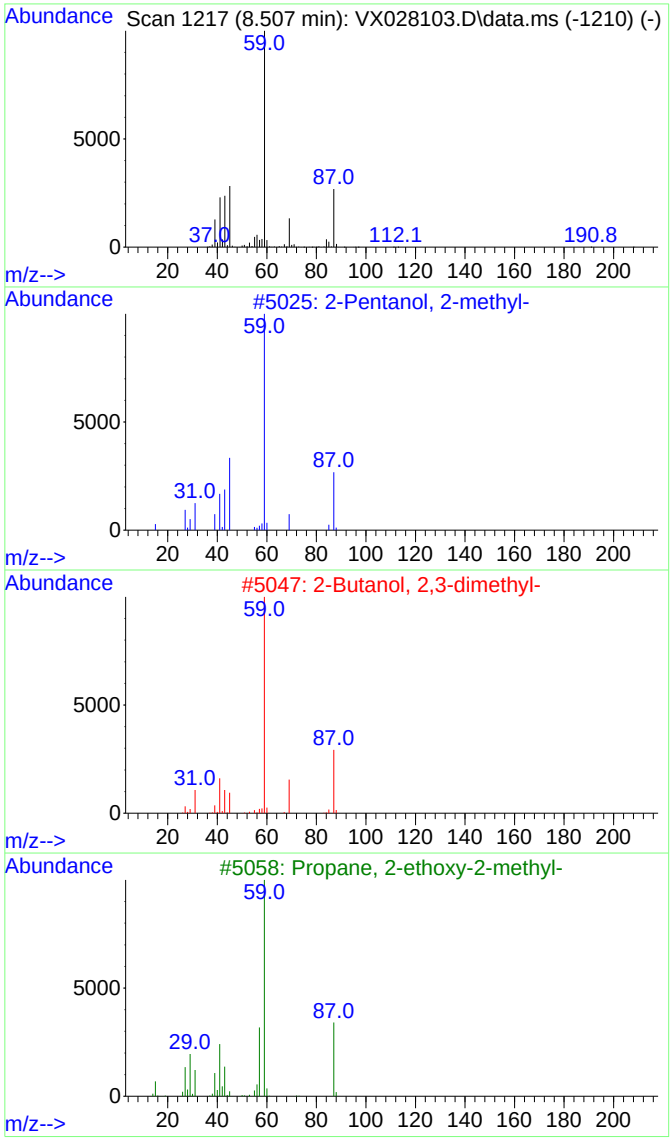
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

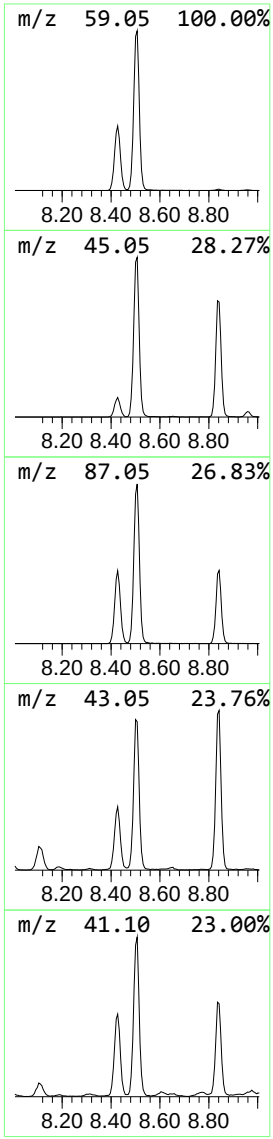
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 2-Pentanol, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.507	75.97 ug/l	1777500	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanol, 2-methyl-	102	C6H14O	000590-36-3	83
2	2-Butanol, 2,3-dimethyl-	102	C6H14O	000594-60-5	64
3	Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	53
4	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	50
5	DL-Lactamide, N,O-dimethyl-	117	C5H11NO2	1000452-56-3	42





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Data File : VX028103.D
Acq On : 14 Apr 2022 13:41
Operator : JC/MD
Sample : N2403-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-29R

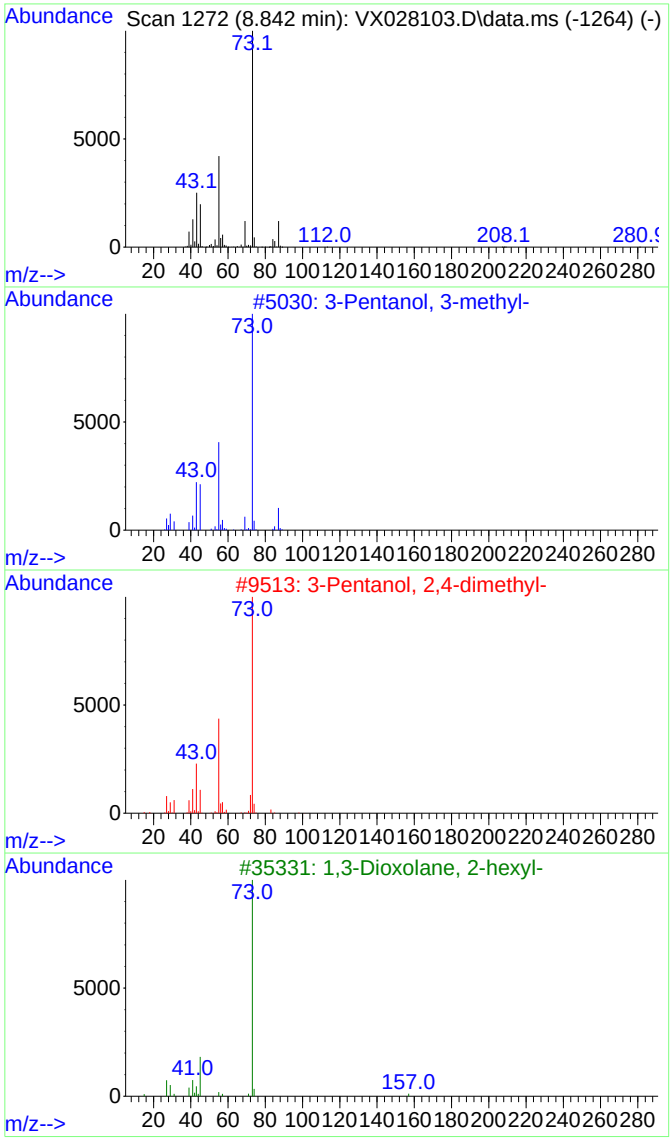
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

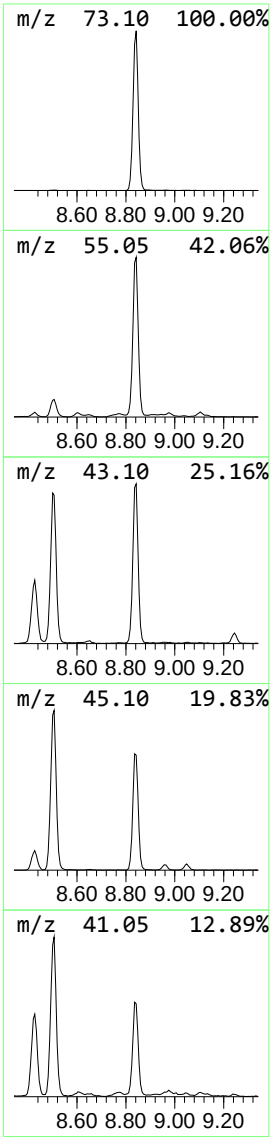
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 3-Pentanol, 3-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.842	73.71 ug/l	1724710	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	83
2	3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	47
3	1,3-Dioxolane, 2-hexyl-	158	C9H18O2	001708-34-5	37
4	Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	25
5	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	16





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028103.D
Acq On : 14 Apr 2022 13:41
Operator : JC/MD
Sample : N2403-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-29R

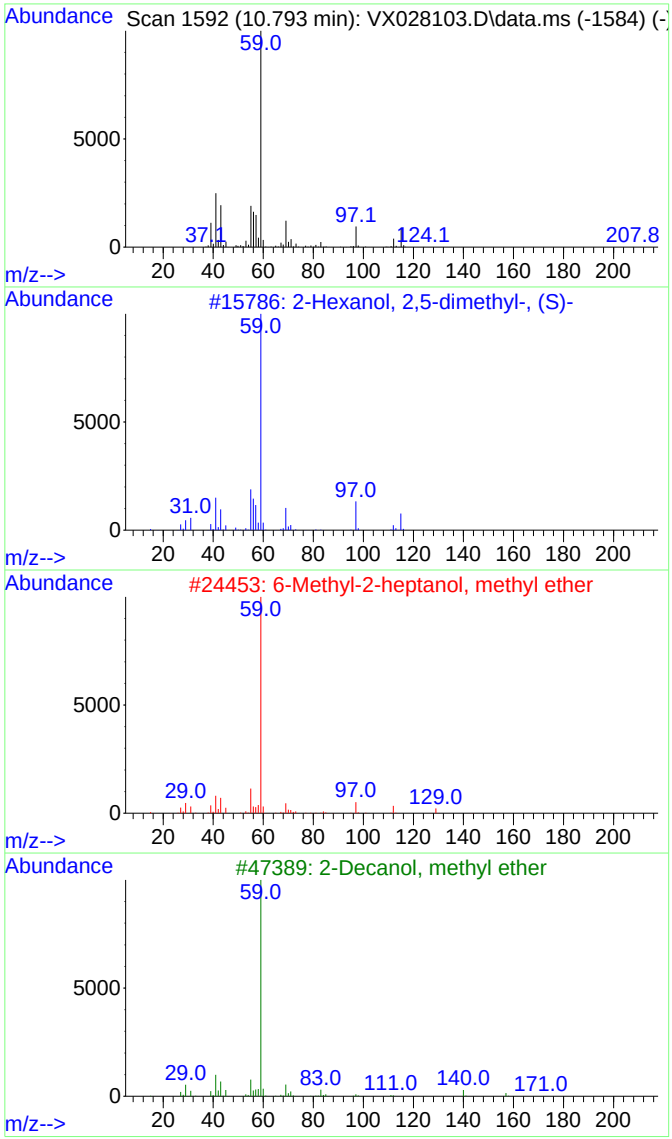
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

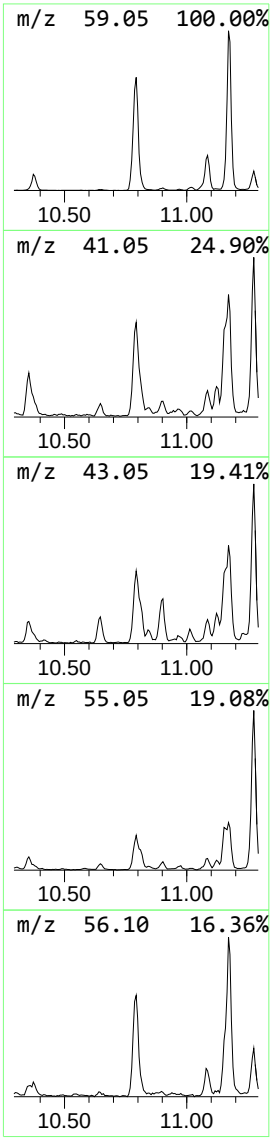
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 2-Hexanol, 2,5-dimethyl-, (S)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.793	18.99 ug/l	444309	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanol, 2,5-dimethyl-, (S)-	130	C8H18O	003730-60-7	90
2	6-Methyl-2-heptanol, methyl ether	144	C9H20O	1000365-52-0	45
3	2-Decanol, methyl ether	172	C11H24O	1000333-83-9	40
4	2-Hydroxy-2,4-dimethyl-3-pentanone	130	C7H14O2	003212-67-7	36
5	3-Nonanol	144	C9H20O	000624-51-1	36





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028103.D
Acq On : 14 Apr 2022 13:41
Operator : JC/MD
Sample : N2403-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-29R

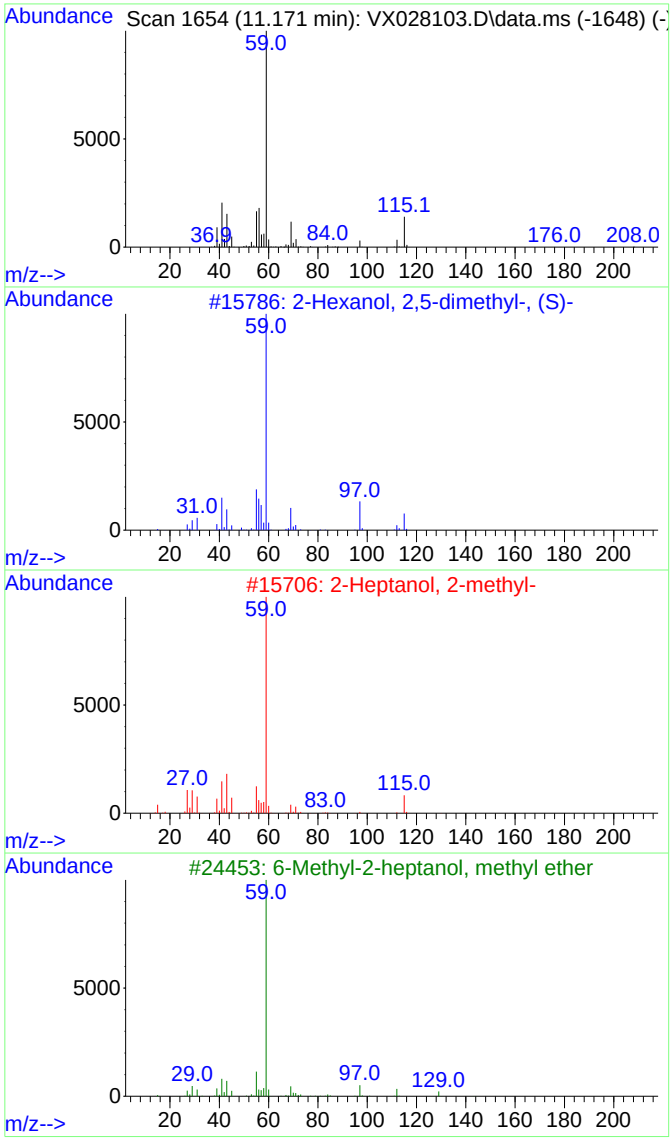
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

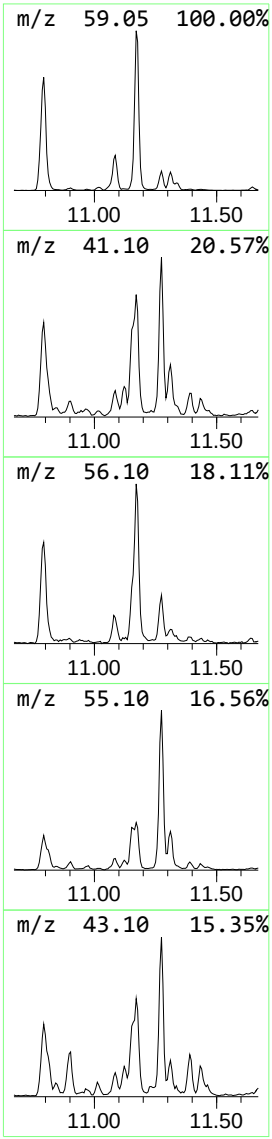
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 2-Heptanol, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.171	27.81 ug/l	662699	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanol, 2,5-dimethyl-, (S)-	130	C8H18O	003730-60-7	78
2	2-Heptanol, 2-methyl-	130	C8H18O	000625-25-2	72
3	6-Methyl-2-heptanol, methyl ether	144	C9H20O	1000365-52-0	59
4	2-Hexanol, 2,3-dimethyl-	130	C8H18O	019550-03-9	50
5	2-Decanol, methyl ether	172	C11H24O	1000333-83-9	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028103.D
Acq On : 14 Apr 2022 13:41
Operator : JC/MD
Sample : N2403-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-29R

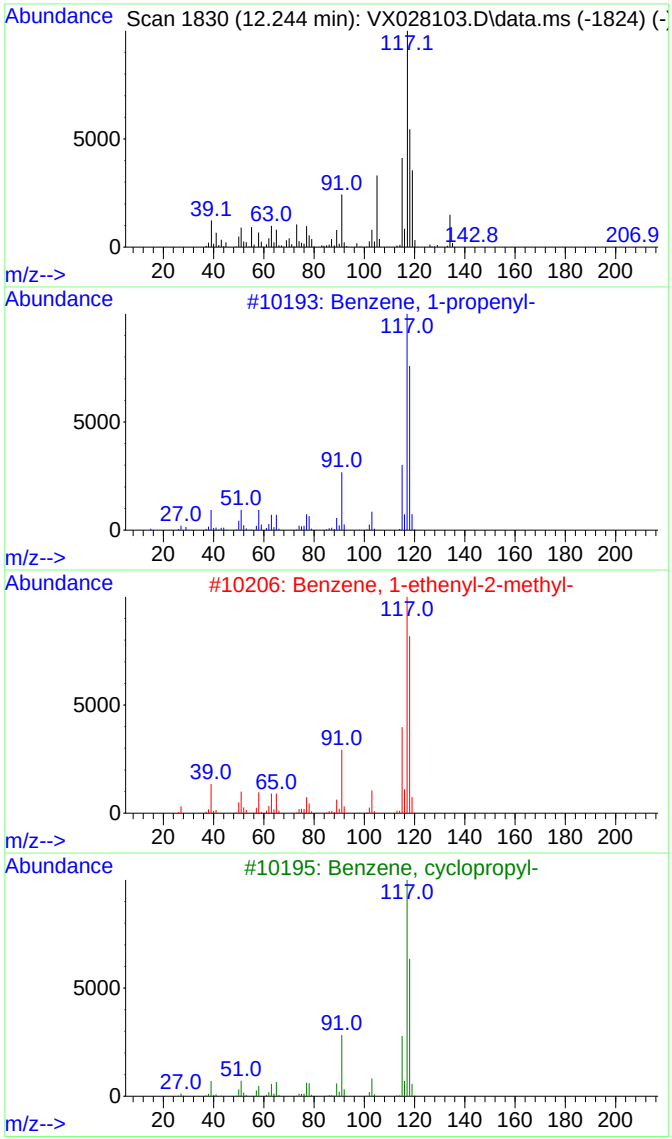
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

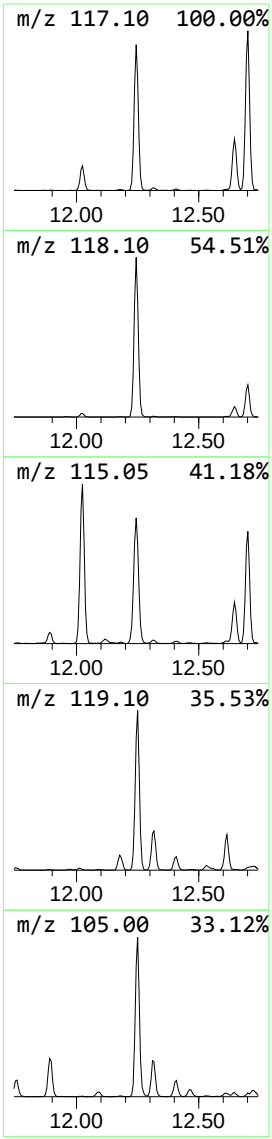
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1-propenyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	54.37 ug/l	1295860	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-propenyl-	118	C9H10	000637-50-3	87
2	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	80
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	55
4	Indane	118	C9H10	000496-11-7	55
5	Benzene, 2-propenyl-	118	C9H10	000300-57-2	55





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028103.D
Acq On : 14 Apr 2022 13:41
Operator : JC/MD
Sample : N2403-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-29R

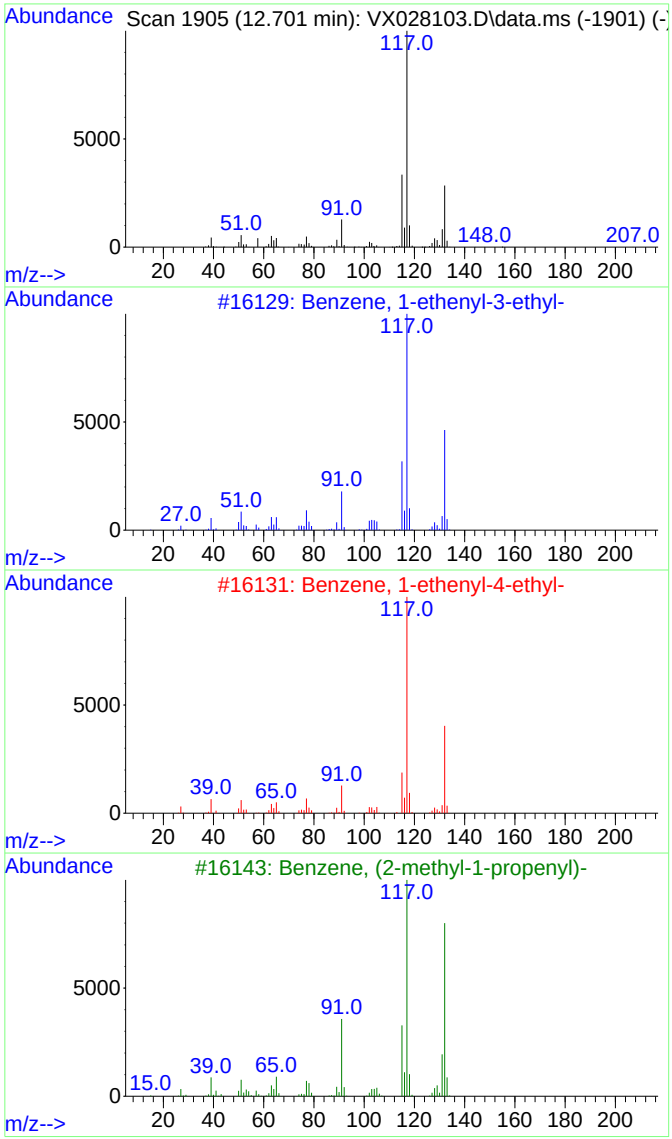
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041222W.M
Quant Title : SW846 8260

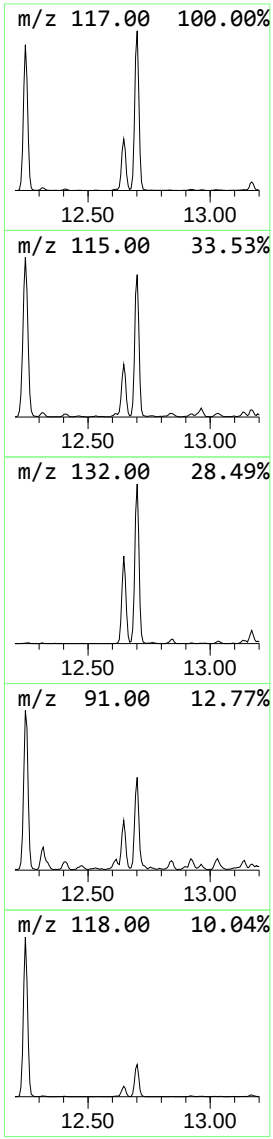
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	29.69 ug/l	707542	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	90
3	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90
4	Indan, 1-methyl-	132	C10H12	000767-58-8	90
5	1-Phenyl-1-butene	132	C10H12	000824-90-8	87





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041422\
Data File : VX028103.D
Acq On : 14 Apr 2022 13:41
Operator : JC/MD
Sample : N2403-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-29R

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-methyl-	1.752	24.5	ug/l	316239	1	5.556	646366	50.0
Butane, 2,3-dim...	2.782	25.4	ug/l	329017	1	5.556	646366	50.0
Amylene hydrate	6.239	201.0	ug/l	3639200	2	6.763	905205	50.0
2-Butanol, 2,3-...	8.427	32.2	ug/l	753862	3	10.055	1169930	50.0
2-Pentanol, 2-m...	8.507	76.0	ug/l	1777500	3	10.055	1169930	50.0
3-Pentanol, 3-m...	8.842	73.7	ug/l	1724710	3	10.055	1169930	50.0
2-Hexanol, 2,5-...	10.793	19.0	ug/l	444309	3	10.055	1169930	50.0
2-Heptanol, 2-m...	11.171	27.8	ug/l	662699	4	12.024	1191650	50.0
Benzene, 1-prop...	12.244	54.4	ug/l	1295860	4	12.024	1191650	50.0
Benzene, 1-ethe...	12.701	29.7	ug/l	707542	4	12.024	1191650	50.0