LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062222\
Data File : VX029707.D
Acq On : 22 Jun 2022 18:20
Operator : JC/MD
Sample : N3202-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-30RE

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: VX029707.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.258	23	28	42	rBV	46719	116189	8.39%	0.841%
2	1.416	50	54	58	rBV	40375	45764	3.31%	0.331%
3	2.130	165	171	177	rVB	16517	29099	2.10%	0.211%
4	2.337	198	205	217	rVB2	22827	49634	3.59%	0.359%
5	2.782	269	278	288	rBV	35838	80643	5.83%	0.584%
6	2.959	300	307	321	rBV	58150	129912	9.38%	0.940%
7	3.111	322	332	344	rVB2	88284	210709	15.22%	1.525%
8	3.764	427	439	445	rBV6	9397	29848	2.16%	0.216%
9	3.843	445	452	459	rVB4	7500	19918	1.44%	0.144%
10	4.434	538	549	563	rVB6	16093	55464	4.01%	0.401%
11	4.898	612	625	632	rBV3	4522	15051	1.09%	0.109%
12	5.196	664	674	689	rVB	21751	72327	5.22%	0.523%
13	5.391	692	706	722	rBV2	150142	434180	31.37%	3.142%
14	5.556	722	733	744	rBV	270925	790916	57.14%	5.724%
15	5.958	789	799	809	rBV	155726	412557	29.80%	2.986%
16	6.233	833	844	858	rBV	497773	1384261	100.00%	10.018%
17	6.367	862	866	874	rVB7	6166	16208	1.17%	0.117%
18	6.763	921	931	942	rBV	414505	1021585	73.80%	7.393%
19	7.178	986	999	1006	rVB5	13868	38476	2.78%	0.278%
20	7.367	1023	1030	1038	rBV5	8544	20671	1.49%	0.150%
21	7.775	1092	1097	1104	rVB3	7802	14234	1.03%	0.103%
22	7.988	1122	1132	1142	rVB4	11114	27859	2.01%	0.202%
23	8.110	1145	1152	1153	rBV	31302	46988	3.39%	0.340%
24	8.147	1153	1158	1163	rVV	110730	197179	14.24%	1.427%
25	8.196	1163	1166	1174	rVB2	8641	14050	1.01%	0.102%
26	8.318	1178	1186	1195	rBV5	7487	19566	1.41%	0.142%
27	8.427	1195	1204	1210	rBV	167132	289718	20.93%	2.097%
28	8.501	1210	1216	1230	rVB	360495	594176	42.92%	4.300%
29	8.653	1230	1241	1250	rBV	842601	1336258	96.53%	9.670%
30	8.842	1265	1272	1281	rBV	400057	645237	46.61%	4.670%
31	8.927	1281	1286	1290	rVV	19444	32276	2.33%	0.234%
32	9.165	1319	1325	1334	rVB	50838	83751	6.05%	0.606%
33	9.446	1364	1371	1377	rBV2	46251	81401	5.88%	0.589%
34	9.506	1377	1381	1385	rVV6	8137	17258	1.25%	0.125%
35	9.561	1385	1390	1395	rVV	92445	146278	10.57%	1.059%
36	9.604	1395	1397	1403	rVV3	16580	24115	1.74%	0.175%

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Instrument :
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ClientSampleId :
MW-30RE

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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Title : SW846 8260

37	9.763	1418	1423	1430	rVV4	8944	14068	1.02%	0.102%
38	10.025	1460	1466	1467	rBV	95722	139150	10.05%	1.007%
39	10.055	1467	1471	1476	rVV	866449	1144015	82.64%	8.279%
40	10.104	1476	1479	1482	rVV	59541	83329	6.02%	0.603%
41	10.147	1482	1486	1493	rVV	163582	226514	16.36%	1.639%
42	10.244	1496	1502	1509	rVV	71289	145097	10.48%	1.050%
43	10.305	1509	1512	1516	rVB	13187	17318	1.25%	0.125%
44	10.354	1516	1520	1526	rBV	16661	23140	1.67%	0.167%
45	10.421	1526	1531	1535	rBV	23186	33190	2.40%	0.240%
46	10.647	1564	1568	1573	rBV2	14750	20200	1.46%	0.146%
47	10.793	1586	1592	1598	rBV2	28781	62048	4.48%	0.449%
48	10.848	1598	1601	1605	rVV2	13345	19798	1.43%	0.143%
49	10.903	1605	1610	1613	rVV	24391	37062	2.68%	0.268%
50	10.939	1613	1616	1619	rVV	18991	27675	2.00%	0.200%
51	11.085	1634	1640	1645	rBV	641553	837760	60.52%	6.063%
52	11.159	1649	1652	1657	rVB2	15807	22204	1.60%	0.161%
53	11.274	1666	1671	1674	rBV	37558	50112	3.62%	0.363%
54	11.311	1674	1677	1686	rVB4	17165	27375	1.98%	0.198%
55	11.390	1686	1690	1695	rBV3	17027	29946	2.16%	0.217%
56	11.451	1695	1700	1709	rVB3	21259	52227	3.77%	0.378%
57	11.616	1721	1727	1735	rBV	15402	25645	1.85%	0.186%
58	11.695	1735	1740	1746	rVV6	13088	28908	2.09%	0.209%
59	11.756	1746	1750	1756	rVB	39520	50598	3.66%	0.366%
60	11.890	1763	1772	1778	rBV3	22080	51596	3.73%	0.373%
61	12.024	1788	1794	1801	rBV	862230	1050050	75.86%	7.599%
62	12.091	1801	1805	1815	rVV2	21996	50339	3.64%	0.364%
63	12.177	1815	1819	1824	rVB	11325	16790	1.21%	0.122%
64	12.244	1824	1830	1836	rBV2	49369	68068	4.92%	0.493%
65	12.317	1837	1842	1852	rVB	64844	84621	6.11%	0.612%
66	12.415	1852	1858	1862	rBV2	25826	35094	2.54%	0.254%
67	12.646	1893	1896	1900	rVV	60398	71057	5.13%	0.514%
68	12.701	1900	1905	1911	rVV	243108	296489	21.42%	2.146%
69	12.841	1921	1928	1933	rBV2	19046	27072	1.96%	0.196%
70	12.927	1934	1942	1945	rBV	35699	50457	3.65%	0.365%
71	13.036	1955	1960	1964	rBV	15537	19351	1.40%	0.140%
72	13.140	1968	1977	1980	rBV2	15469	28038	2.03%	0.203%
73	13.195	1984	1986	1991	rVB	17874	19889	1.44%	0.144%
74	13.292	1998	2002	2006	rVB2	12087	15968	1.15%	0.116%
75	13.347	2006	2011	2016	rVV2	19760	27188	1.96%	0.197%
76	13.427	2019	2024	2031	rVB2	45639	61430	4.44%	0.445%

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ALS Vial : 24 Sample Multiplier: 1

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MSVOA_X
ClientSampleId :
MW-30RE

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	13.542	2039	2043	2046	rBV	15491	20782	1.50%	0.150%
78	13.640	2053	2059	2068	rVB2	22158	45819	3.31%	0.332%
79	13.780	2074	2082	2090	rBV2	13608	25361	1.83%	0.184%
80	13.926	2100	2106	2111	rBV2	22406	27929	2.02%	0.202%
81	14.597	2210	2216	2220	rBV	15628	23409	1.69%	0.169%
82	14.640	2220	2223	2228	rVB	14349	18609	1.34%	0.135%
83	14.786	2242	2247	2252	rBV3	13513	21442	1.55%	0.155%

Sum of corrected areas: 13817983

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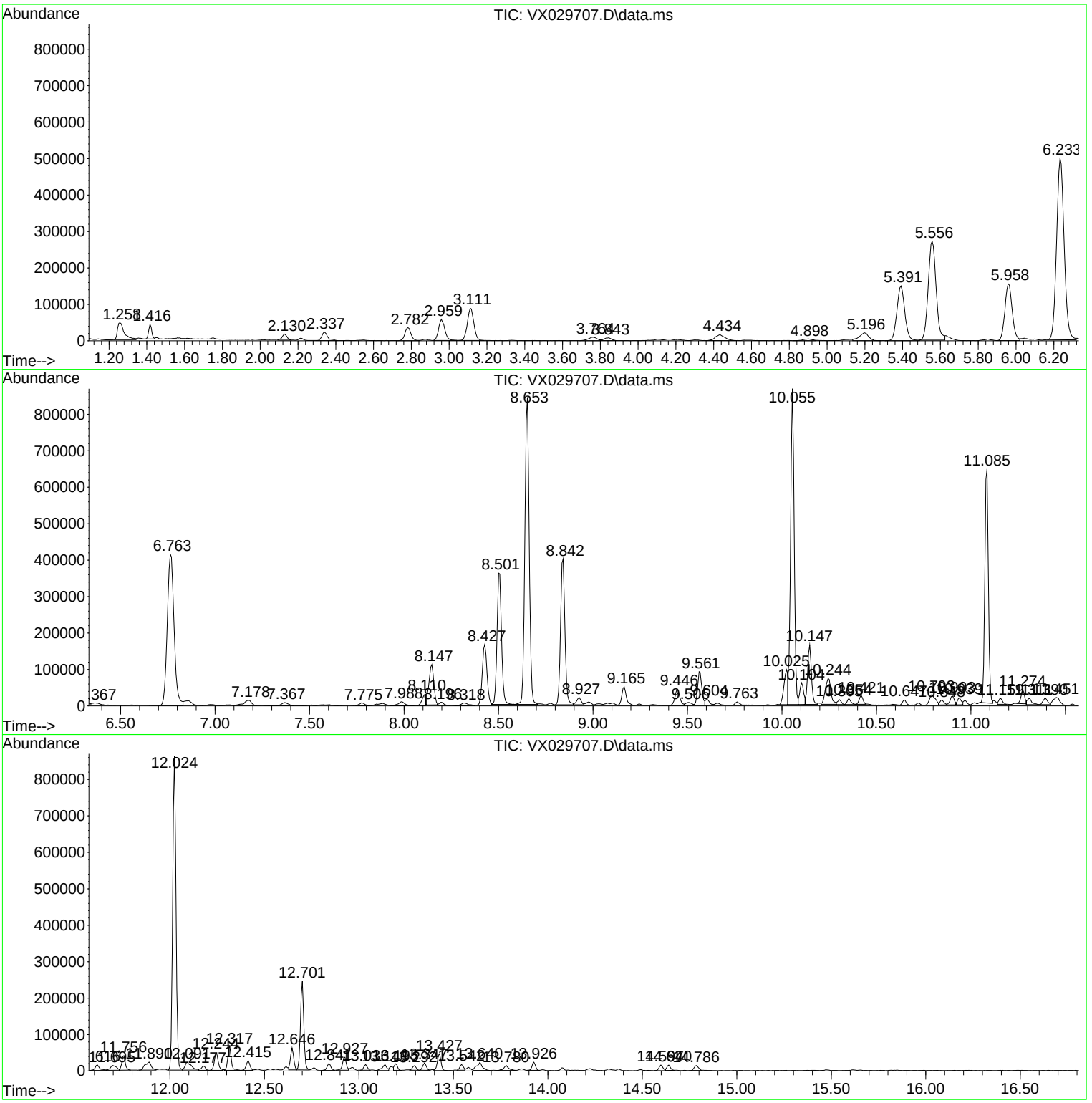
Instrument :
MSVOA_X
ClientSampleId :
MW-30RE

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



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Data File : VX029707.D
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Operator : JC/MD
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ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-30RE

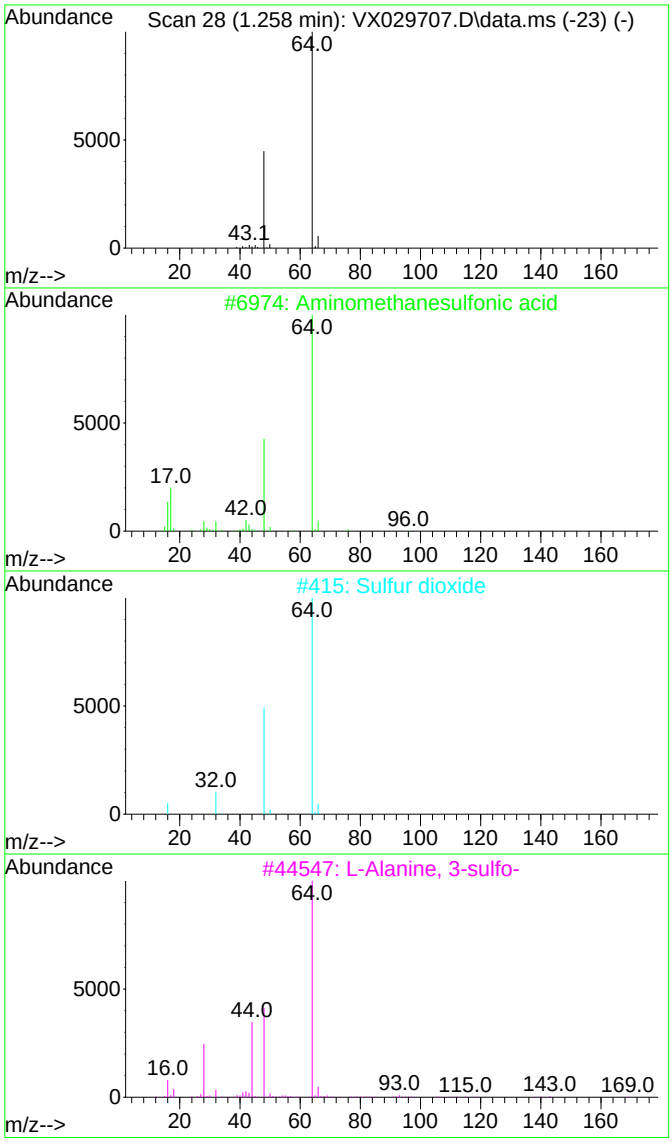
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

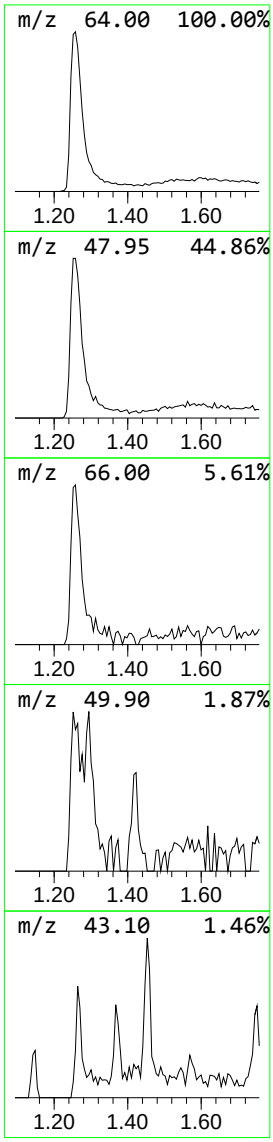
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Aminomethanesulfonic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.258	7.35 ug/l	116189	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	83
2	Sulfur dioxide	64	O2S	007446-09-5	83
3	L-Alanine, 3-sulfo-	169	C3H7NO5S	000498-40-8	9
4	Ethene, 1,1-difluoro-	64	C2H2F2	000075-38-7	5
5	Ethene, 1,2-difluoro-	64	C2H2F2	001691-13-0	4





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ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-30RE

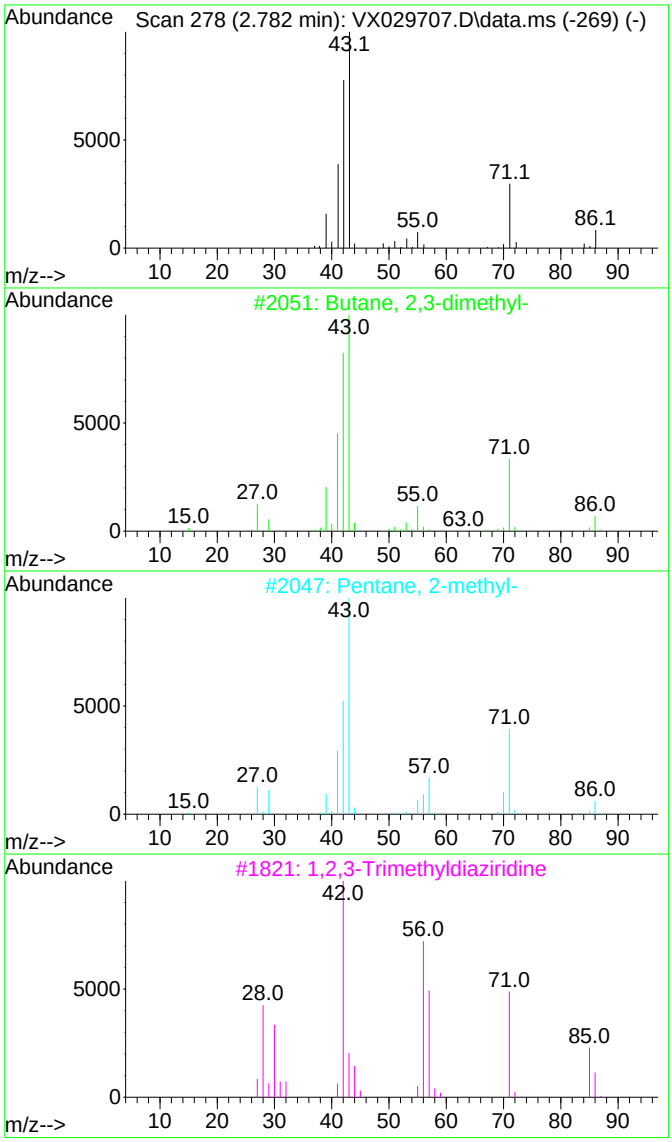
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

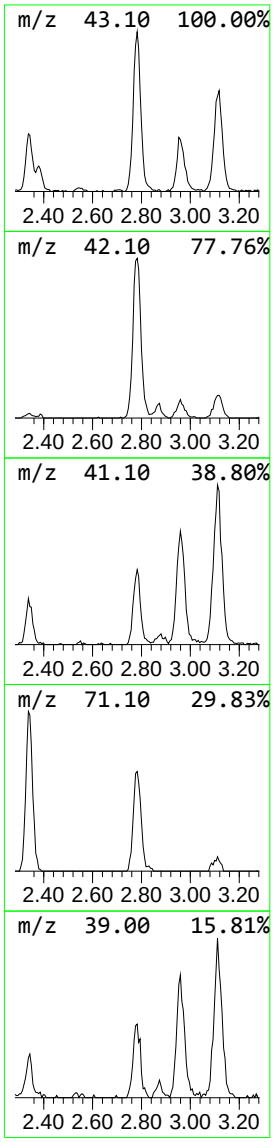
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Butane, 2,3-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.782	5.10 ug/l	80643	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	64
3	1,2,3-Trimethyldiaziridine	86	C4H10N2	113604-56-1	36
4	Tetrahydrofuran	72	C4H8O	000109-99-9	36
5	Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	9





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Data File : VX029707.D
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Operator : JC/MD
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-30RE

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

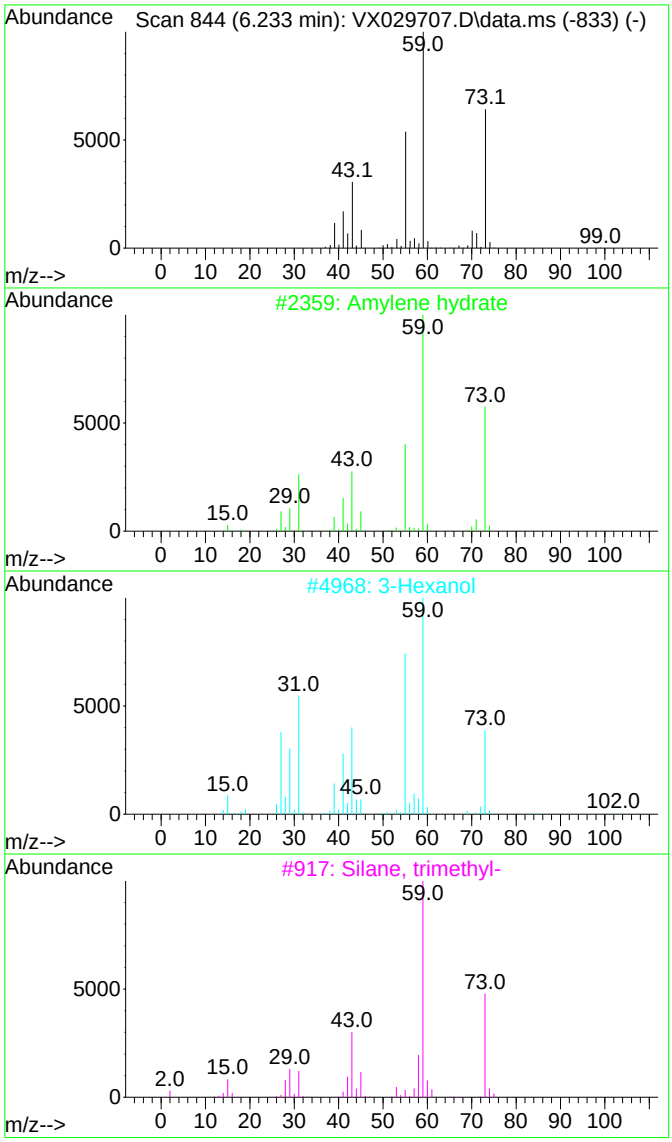
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

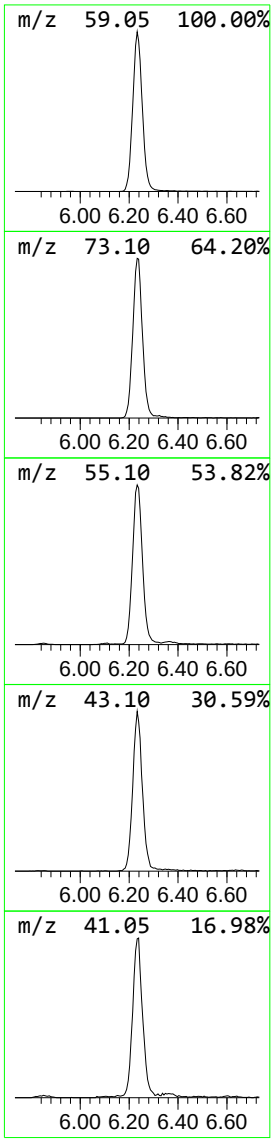
****

Peak Number 3 Amylene hydrate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.233	67.75 ug/l	1384260	1,4-Difluorobenzene	6.769

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	40
3	Silane, trimethyl-	74	C3H10Si	000993-07-7	39
4	3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	38
5	1-Hepten-4-ol	114	C7H14O	003521-91-3	9





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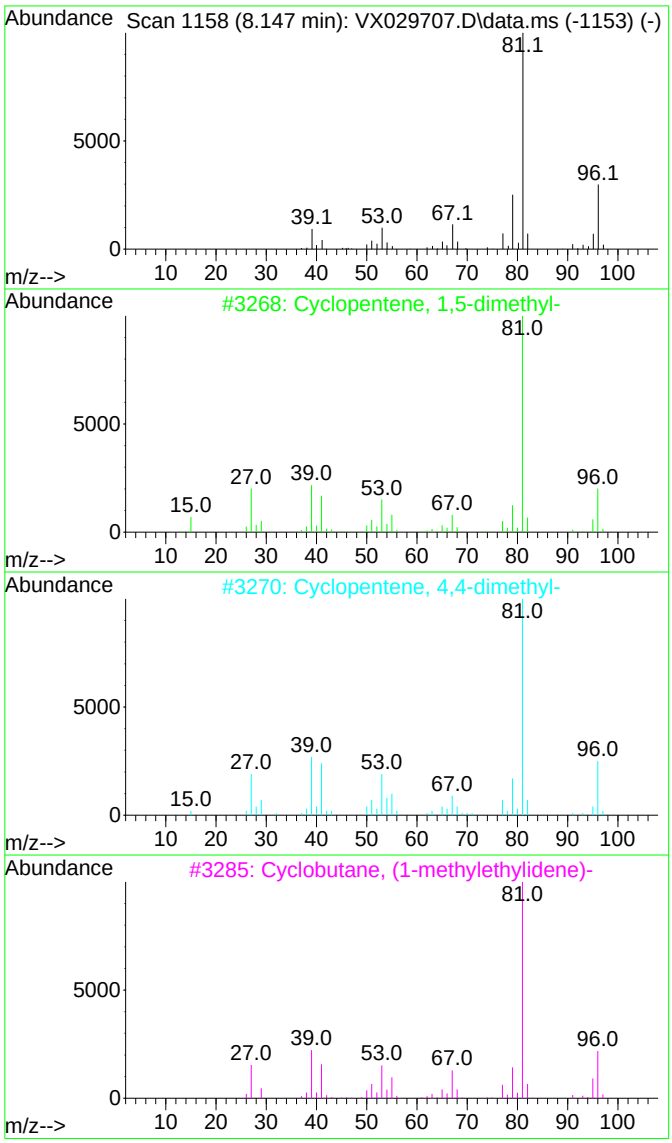
Quant Title : SW846 8260

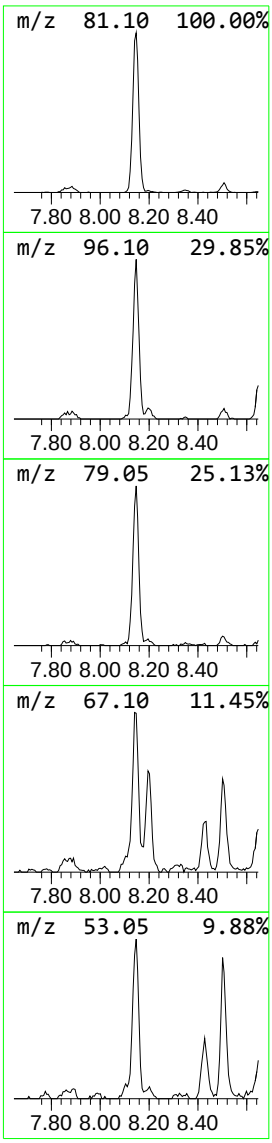
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Cyclopentene, 1,5-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.147	9.65 ug/l	197179	1,4-Difluorobenzene	6.769

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentene, 1,5-dimethyl-	96	C7H12	016491-15-9	86
2	Cyclopentene, 4,4-dimethyl-	96	C7H12	019037-72-0	86
3	Cyclobutane, (1-methylethylidene)-	96	C7H12	001528-22-9	72
4	1,4-Pentadiene, 3,3-dimethyl-	96	C7H12	001112-35-2	72
5	Cyclopropane, 1-methyl-1-isoprop...	96	C7H12	003422-07-9	72





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ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-30RE

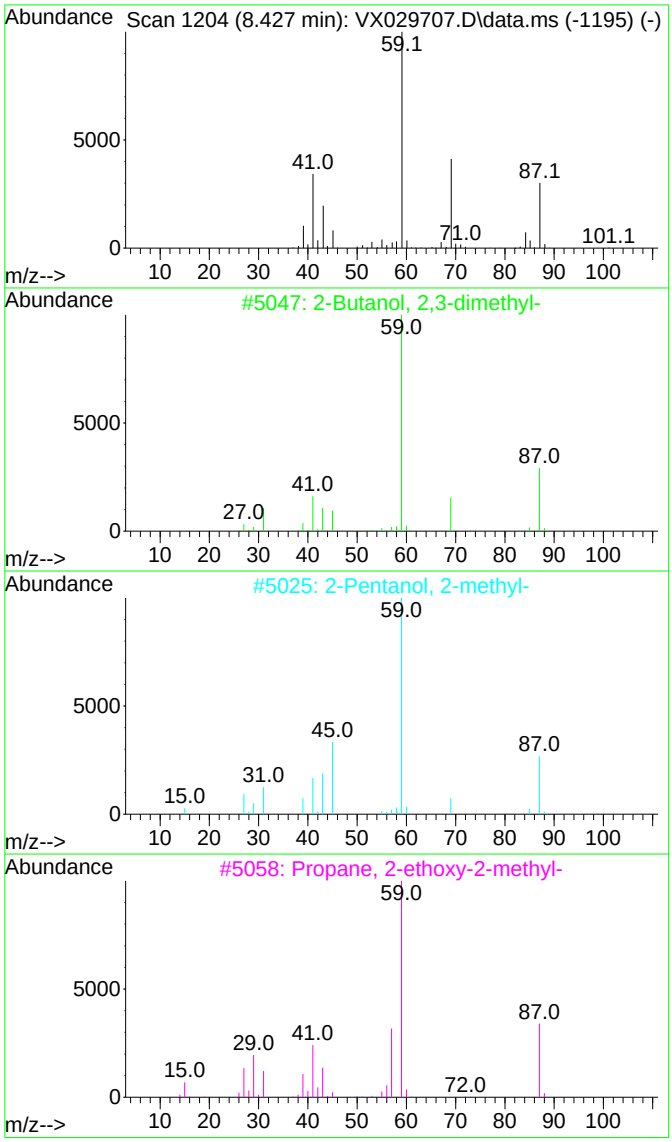
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

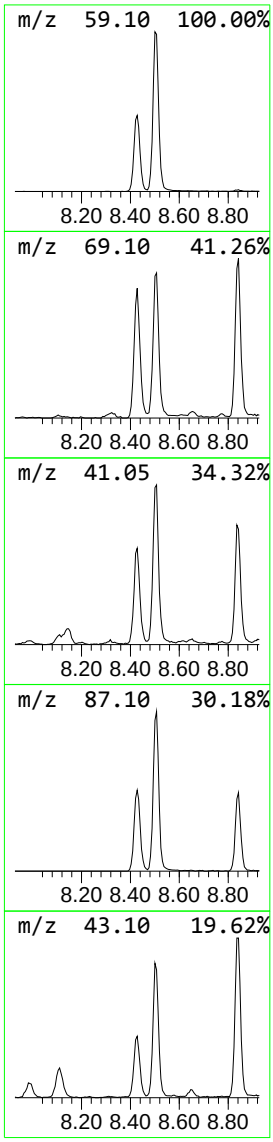
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.427	12.66 ug/l	289718	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanol, 2,3-dimethyl-	102	C6H14O	000594-60-5	78
2	2-Pentanol, 2-methyl-	102	C6H14O	000590-36-3	72
3	Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	56
4	4-Pentene-2-ol, 2-methyl	100	C6H12O	000624-97-5	50
5	3-Pentanol, 2,2-dimethyl-	116	C7H16O	003970-62-5	39





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Data File : VX029707.D
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ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-30RE

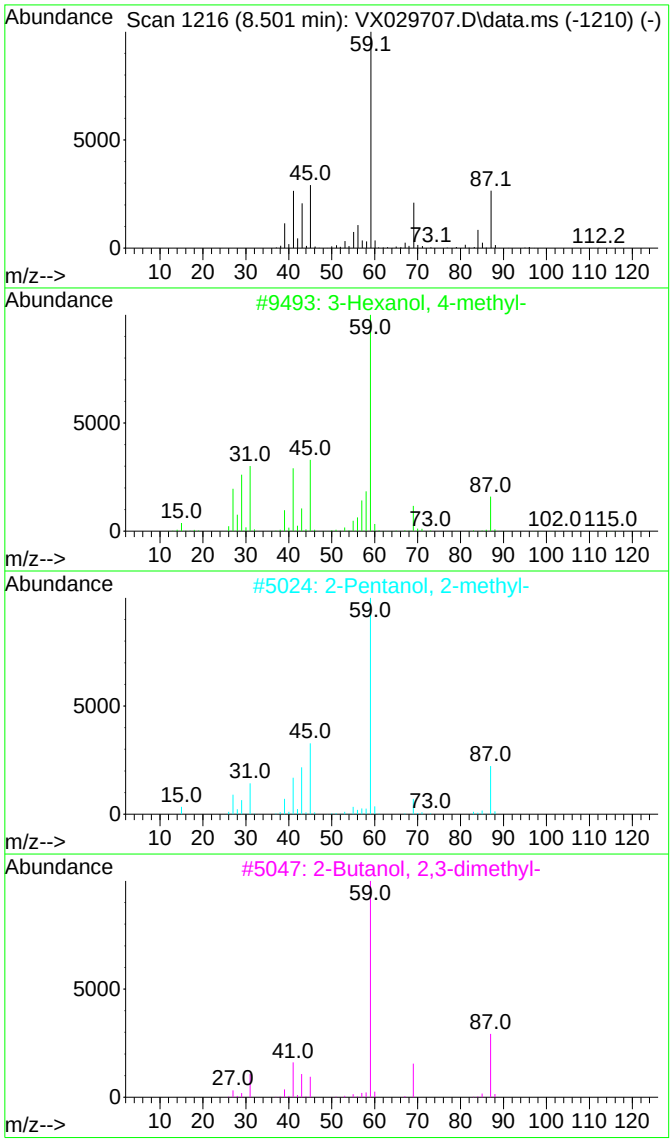
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

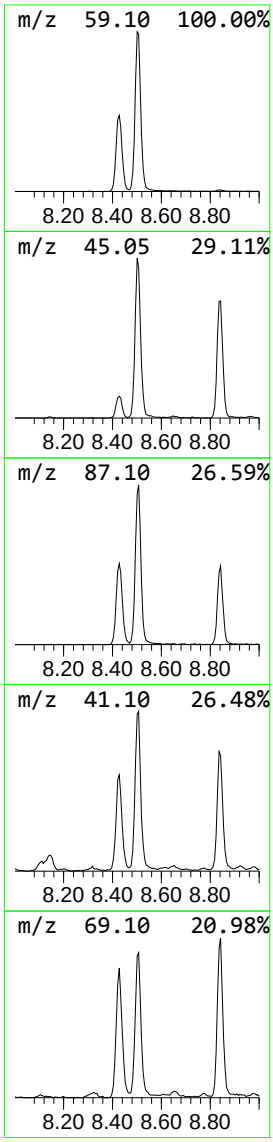
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 3-Hexanol, 4-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.501	25.97 ug/l	594176	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	78
2	2-Pentanol, 2-methyl-	102	C6H14O	000590-36-3	78
3	2-Butanol, 2,3-dimethyl-	102	C6H14O	000594-60-5	64
4	3-Pentanol, 2,2-dimethyl-	116	C7H16O	003970-62-5	56
5	3-Hexanol, 5-methyl-	116	C7H16O	000623-55-2	56





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062222\
Data File : VX029707.D
Acq On : 22 Jun 2022 18:20
Operator : JC/MD
Sample : N3202-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-30RE

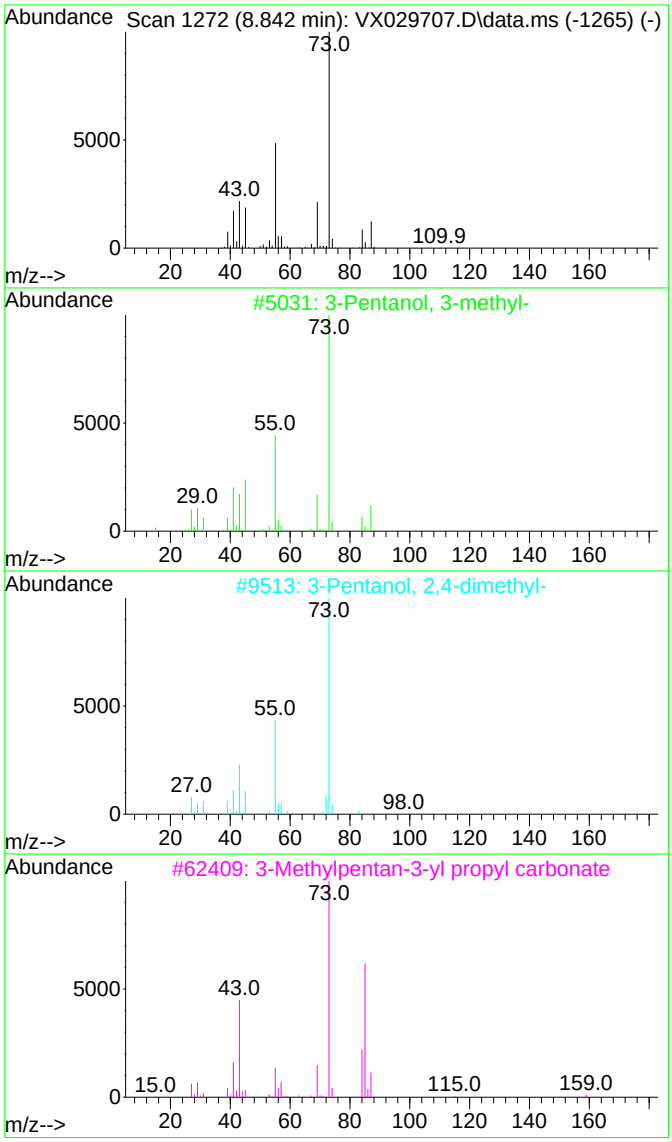
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

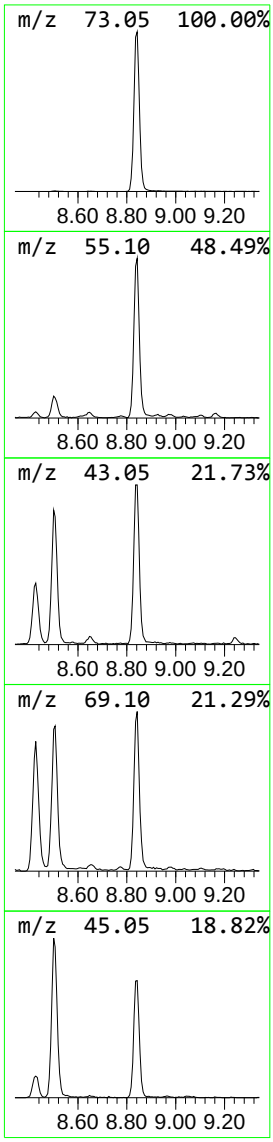
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 3-Pentanol, 3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.842	28.20 ug/l	645237	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	53
3	3-Methylpentan-3-yl propyl carbo...	188	C10H20O3	1000372-82-3	33
4	Methane, isothiocyanato-	73	C2H3NS	000556-61-6	32
5	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	10





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062222\
Data File : VX029707.D
Acq On : 22 Jun 2022 18:20
Operator : JC/MD
Sample : N3202-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-30RE

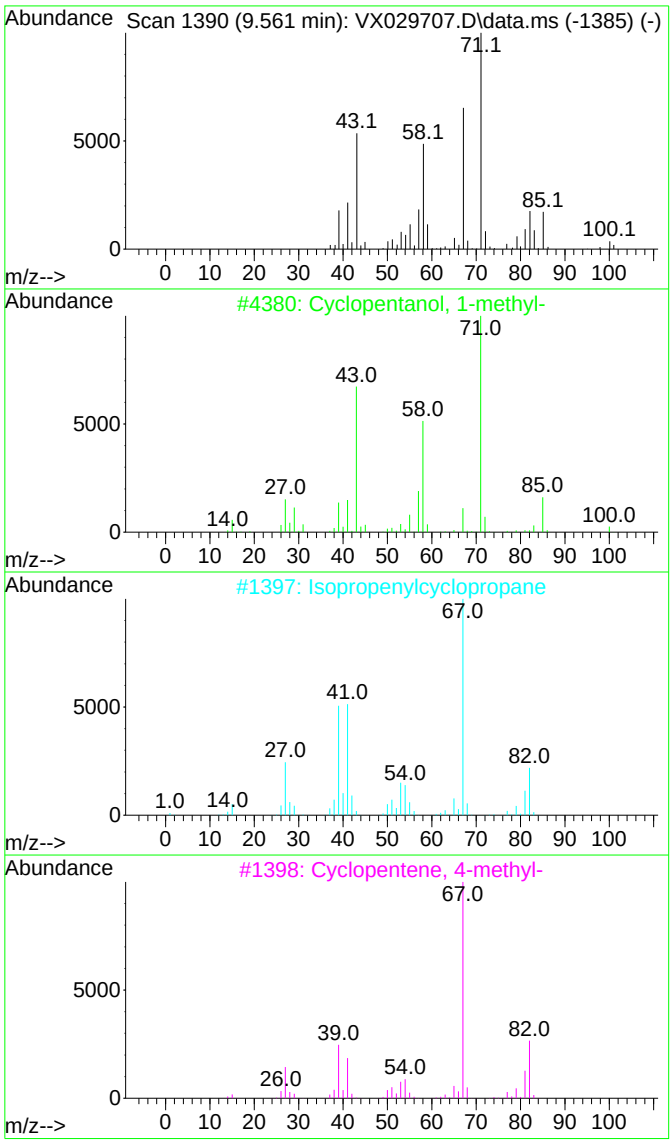
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

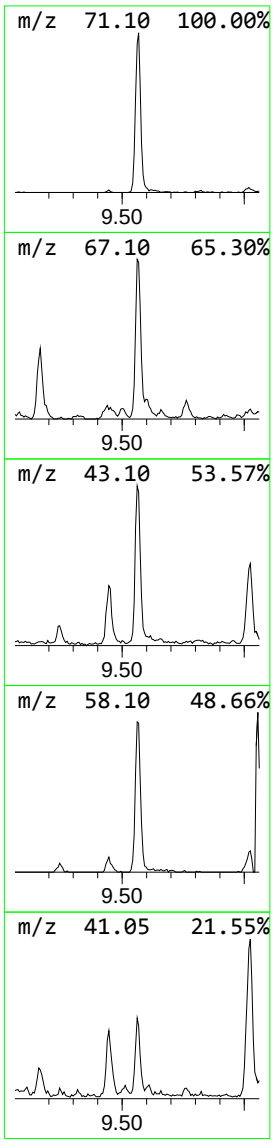
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Cyclopentanol, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.561	6.39 ug/l	146278	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentanol, 1-methyl-	100	C6H12O	001462-03-9	58
2	Isopropenylcyclopropane	82	C6H10	004663-22-3	25
3	Cyclopentene, 4-methyl-	82	C6H10	001759-81-5	25
4	Cyclopentene, 1-methyl-	82	C6H10	000693-89-0	25
5	1,4-Pentadiene, 2-methyl-	82	C6H10	000763-30-4	25





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062222\
Data File : VX029707.D
Acq On : 22 Jun 2022 18:20
Operator : JC/MD
Sample : N3202-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-30RE

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

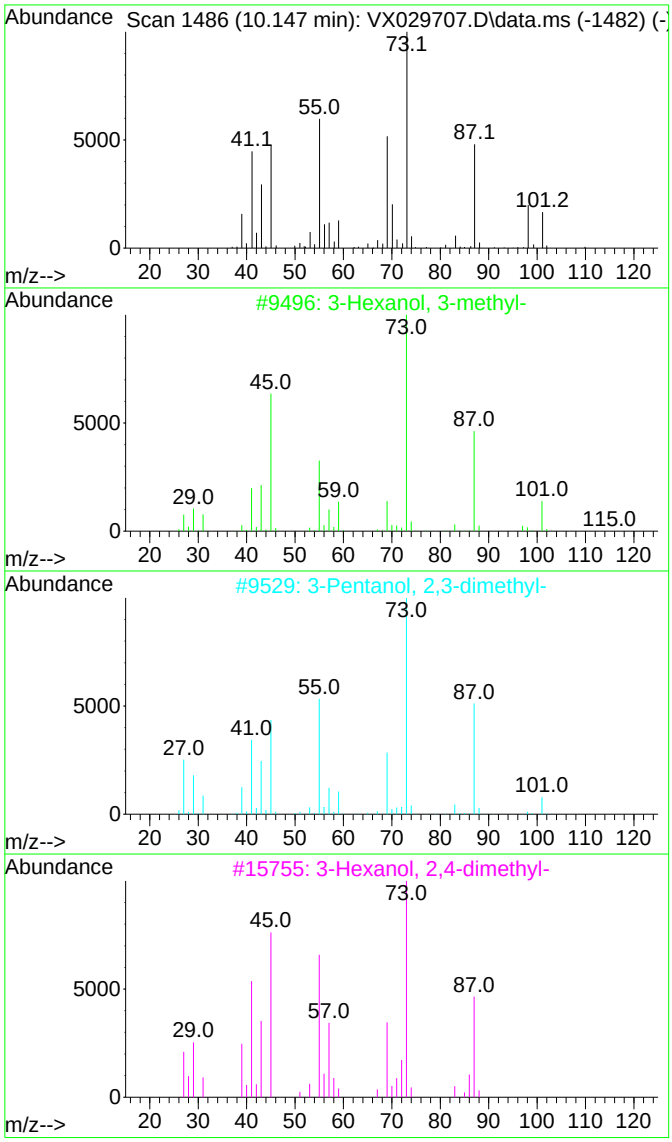
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

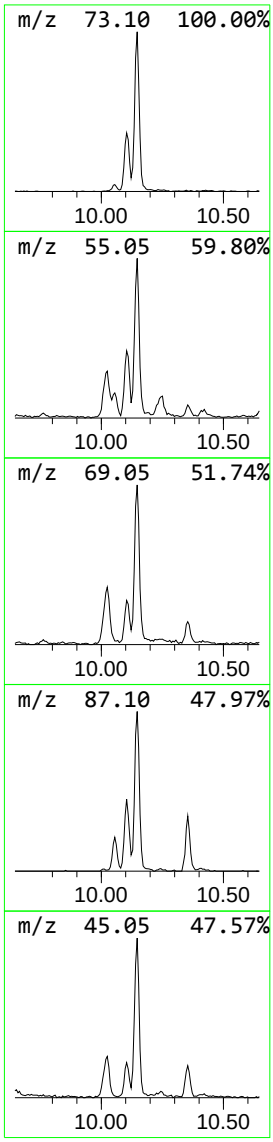
****

Peak Number 9 3-Hexanol, 3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.147	9.90 ug/l	226514	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanol, 3-methyl-	116	C7H16O	000597-96-6	59
2	3-Pentanol, 2,3-dimethyl-	116	C7H16O	000595-41-5	50
3	3-Hexanol, 2,4-dimethyl-	130	C8H18O	013432-25-2	47
4	3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	43
5	4-Heptanol, 2,6-dimethyl-	144	C9H20O	000108-82-7	37





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Data File : VX029707.D
Acq On : 22 Jun 2022 18:20
Operator : JC/MD
Sample : N3202-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-30RE

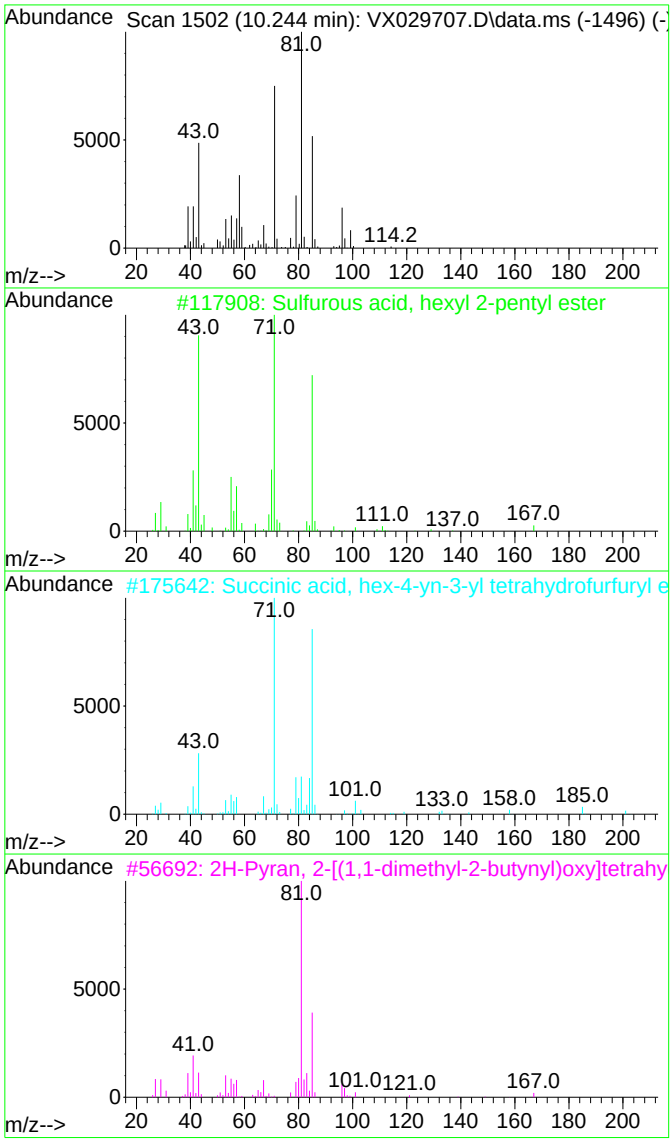
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

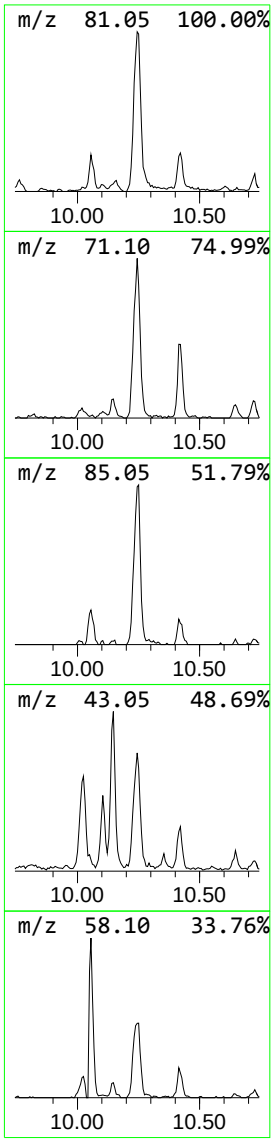
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 unknown10.244 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.244	6.34 ug/l	145097	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Sulfurous acid, hexyl 2-pentyl e...	236	C11H24O3S	1000309-15-6	35
2	Succinic acid, hex-4-yn-3-yl tet...	282	C15H22O5	1000390-71-7	32
3	2H-Pyran, 2-[(1,1-dimethyl-2-but...	182	C11H18O2	024642-03-3	32
4	Cyclobutane, (1-methylethylidene)-	96	C7H12	001528-22-9	25
5	Cyclopentene, 1,5-dimethyl-	96	C7H12	016491-15-9	25





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Data File : VX029707.D
Acq On : 22 Jun 2022 18:20
Operator : JC/MD
Sample : N3202-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-30RE

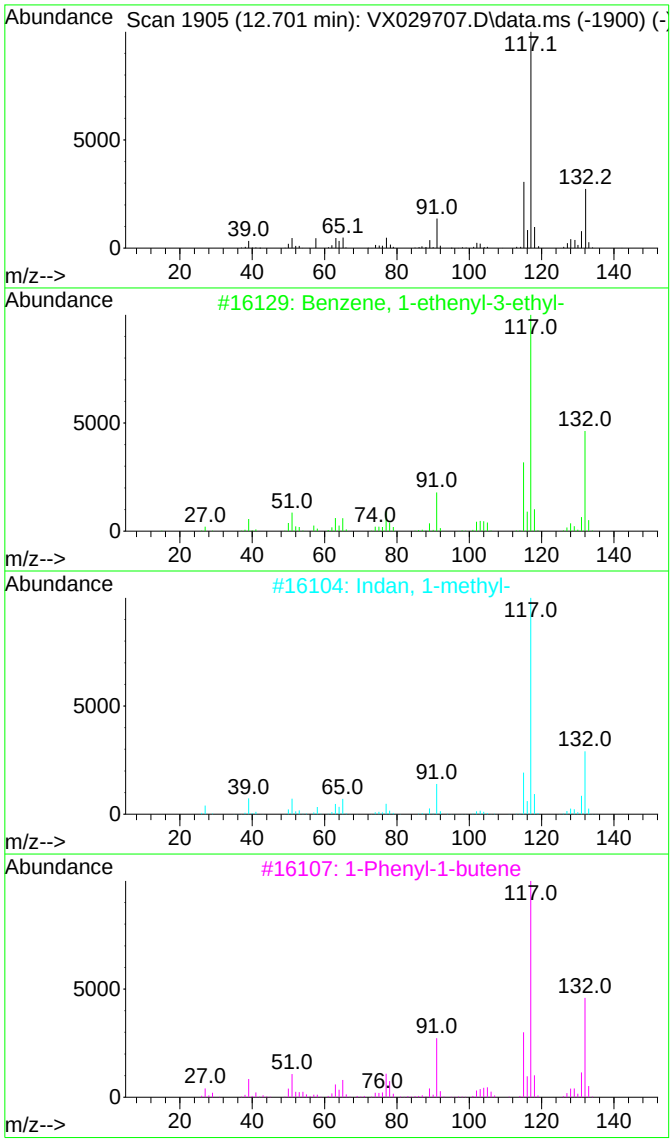
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
Quant Title : SW846 8260

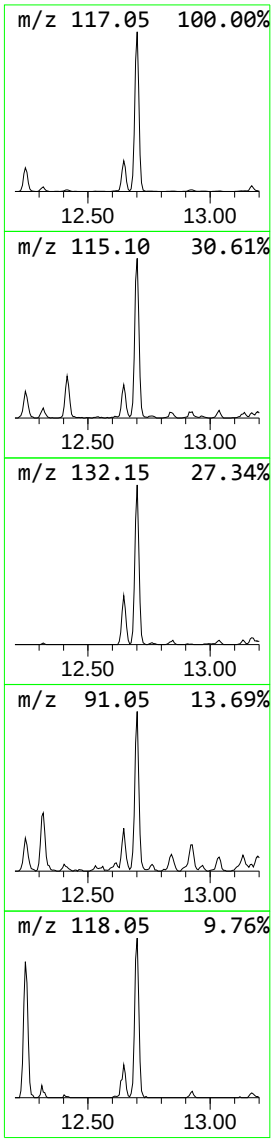
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	14.12 ug/l	296489	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	Indan, 1-methyl-	132	C10H12	000767-58-8	90
3	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
4	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	86





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062222\
Data File : VX029707.D
Acq On : 22 Jun 2022 18:20
Operator : JC/MD
Sample : N3202-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-30RE

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
Aminomethanesul...	1.258	7.3	ug/l	116189	1	5.556	790916	50.0
Butane, 2,3-dim...	2.782	5.1	ug/l	80643	1	5.556	790916	50.0
Amylene hydrate	6.233	67.8	ug/l	1384260	2	6.769	1021590	50.0
Cyclopentene, 1...	8.147	9.7	ug/l	197179	2	6.769	1021590	50.0
2-Butanol, 2,3-...	8.427	12.7	ug/l	289718	3	10.055	1144020	50.0
3-Hexanol, 4-me...	8.501	26.0	ug/l	594176	3	10.055	1144020	50.0
3-Pentanol, 3-m...	8.842	28.2	ug/l	645237	3	10.055	1144020	50.0
Cyclopentanol, ...	9.561	6.4	ug/l	146278	3	10.055	1144020	50.0
3-Hexanol, 3-me...	10.147	9.9	ug/l	226514	3	10.055	1144020	50.0
unknown10.244	10.244	6.3	ug/l	145097	3	10.055	1144020	50.0
Benzene, 1-ethe...	12.701	14.1	ug/l	296489	4	12.024	1050050	50.0