LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX052622\
Data File : VX029011.D
Acq On : 26 May 2022 19:55
Operator : JC/MD
Sample : N3059-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
4922A

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: VX029011.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.252	23	27	42	rBV	599420	1331102	73.73%	4.895%
2	1.453	57	60	72	rVV3	32106	68629	3.80%	0.252%
3	1.544	72	75	82	rVB	39198	48473	2.68%	0.178%
4	1.745	104	108	114	rVB	26365	37518	2.08%	0.138%
5	1.953	138	142	150	rVB2	15655	24835	1.38%	0.091%
6	2.062	155	160	170	rVB4	10286	22841	1.27%	0.084%
7	2.373	205	211	216	rBV	367464	602298	33.36%	2.215%
8	2.422	216	219	230	rVB	67862	107355	5.95%	0.395%
9	2.532	230	237	251	rBV	162757	361442	20.02%	1.329%
10	4.556	560	569	589	rVB	56578	165097	9.14%	0.607%
11	5.007	633	643	657	rBV2	14770	53287	2.95%	0.196%
12	5.385	692	705	718	rBV	216575	633851	35.11%	2.331%
13	5.556	722	733	752	rVB	341067	962527	53.31%	3.539%
14	5.958	785	799	808	rBV	222791	585179	32.41%	2.152%
15	6.037	808	812	822	rVB2	20554	53637	2.97%	0.197%
16	6.763	921	931	944	rBV	545473	1303801	72.21%	4.794%
17	8.580	1221	1229	1234	rBV	71270	117806	6.52%	0.433%
18	8.653	1234	1241	1247	rVV	1137925	1788174	99.04%	6.576%
19	8.720	1247	1252	1259	rVB	289441	449264	24.88%	1.652%
20	9.031	1295	1303	1314	rBV	62096	105865	5.86%	0.389%
21	9.945	1449	1453	1460	rVB	13272	19328	1.07%	0.071%
22	10.055	1465	1471	1484	rVB	1223289	1638798	90.77%	6.026%
23	10.195	1489	1494	1500	rBV	187773	247036	13.68%	0.908%
24	10.299	1506	1511	1518	rBV	799695	1095298	60.67%	4.028%
25	10.506	1538	1545	1553	rBV	36519	51793	2.87%	0.190%
26	10.640	1562	1567	1572	rBV	318015	412842	22.87%	1.518%
27	10.683	1572	1574	1583	rVB	23077	30836	1.71%	0.113%
28	10.774	1583	1589	1594	rBV	25925	39812	2.21%	0.146%
29	10.860	1594	1603	1614	rVB10	9918	28276	1.57%	0.104%
30	11.085	1628	1640	1651	rBV	1001495	1361365	75.40%	5.006%
31	11.305	1669	1676	1679	rBV	21807	41392	2.29%	0.152%
32	11.378	1684	1688	1694	rVV	52537	105997	5.87%	0.390%
33	11.439	1694	1698	1703	rVV4	59365	136137	7.54%	0.501%
34	11.481	1703	1705	1715	rVB4	39331	68504	3.79%	0.252%
35	11.616	1722	1727	1731	rBV	25220	34938	1.94%	0.128%
36	11.756	1745	1750	1758	rVB	174189	230494	12.77%	0.848%

LSC Area Percent Report

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Data File : VX029011.D
Acq On : 26 May 2022 19:55
Operator : JC/MD
Sample : N3059-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
4922A

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	11.835	1758	1763	1766	rBV	114277	156773	8.68%	0.576%
38	11.890	1766	1772	1780	rVV2	646235	1493602	82.73%	5.492%
39	11.951	1780	1782	1789	rVV2	140956	258672	14.33%	0.951%
40	12.024	1789	1794	1801	rVV	1355481	1805483	100.00%	6.639%
41	12.103	1801	1807	1811	rVV2	415795	613434	33.98%	2.256%
42	12.152	1811	1815	1818	rVV2	85012	159231	8.82%	0.586%
43	12.219	1822	1826	1831	rVV	649007	955869	52.94%	3.515%
44	12.280	1831	1836	1839	rVV3	372757	765775	42.41%	2.816%
45	12.311	1839	1841	1847	rVV2	366801	490311	27.16%	1.803%
46	12.384	1847	1853	1857	rVV4	78424	195838	10.85%	0.720%
47	12.426	1857	1860	1863	rVB	89335	103662	5.74%	0.381%
48	12.463	1863	1866	1873	rVB2	28373	43724	2.42%	0.161%
49	12.530	1873	1877	1879	rBV2	23922	32135	1.78%	0.118%
50	12.579	1882	1885	1889	rBV2	75038	84761	4.69%	0.312%
51	12.768	1912	1916	1918	rBV	185032	245244	13.58%	0.902%
52	12.859	1927	1931	1937	rBV	371090	519774	28.79%	1.911%
53	12.969	1945	1949	1951	rBV2	89679	121672	6.74%	0.447%
54	12.993	1951	1953	1956	rVV2	110488	150279	8.32%	0.553%
55	13.024	1956	1958	1960	rVV	63503	84197	4.66%	0.310%
56	13.067	1960	1965	1968	rVV3	123307	289281	16.02%	1.064%
57	13.103	1968	1971	1974	rVV	304386	479210	26.54%	1.762%
58	13.140	1974	1977	1986	rVV5	252955	552669	30.61%	2.032%
59	13.219	1986	1990	1994	rVV	80107	123661	6.85%	0.455%
60	13.268	1994	1998	2000	rVV	148690	188361	10.43%	0.693%
61	13.292	2000	2002	2007	rVB3	95704	113281	6.27%	0.417%
62	13.384	2014	2017	2020	rBV2	29121	33312	1.85%	0.122%
63	13.420	2020	2023	2032	rVB	103609	166331	9.21%	0.612%
64	13.512	2033	2038	2046	rBV2	97885	168312	9.32%	0.619%
65	13.585	2047	2050	2058	rVB8	30309	48507	2.69%	0.178%
66	13.682	2060	2066	2068	rBV5	31229	64223	3.56%	0.236%
67	13.737	2071	2075	2078	rBV4	60270	84348	4.67%	0.310%
68	13.774	2078	2081	2088	rVB2	94968	140529	7.78%	0.517%
69	13.847	2088	2093	2098	rBV3	69153	151146	8.37%	0.556%
70	13.890	2098	2100	2103	rBV3	21596	26398	1.46%	0.097%
71	14.042	2120	2125	2128	rBV	168342	189982	10.52%	0.699%
72	14.072	2128	2130	2133	rBV2	18428	21607	1.20%	0.079%
73	14.231	2151	2156	2161	rVB	88954	145216	8.04%	0.534%
74	14.322	2166	2171	2175	rBV6	15003	28736	1.59%	0.106%
75	14.371	2176	2179	2181	rBV2	19389	23492	1.30%	0.086%
76	14.408	2181	2185	2188	rBV3	22225	33387	1.85%	0.123%

LSC Area Percent Report

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Operator : JC/MD
Sample : N3059-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
4922A

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	14.481	2193	2197	2199	rBV3	66432	100992	5.59%	0.371%
78	14.566	2207	2211	2215	rVB2	29688	40984	2.27%	0.151%
79	14.615	2215	2219	2221	rBV2	88195	109077	6.04%	0.401%
80	14.670	2226	2228	2234	rVB3	29554	41192	2.28%	0.151%
81	14.755	2239	2242	2245	rBV	177096	188611	10.45%	0.694%
82	14.786	2245	2247	2251	rVB4	26315	32572	1.80%	0.120%
83	15.042	2285	2289	2293	rBV6	9823	18245	1.01%	0.067%
84	15.103	2295	2299	2302	rBV4	17954	32133	1.78%	0.118%
85	15.182	2308	2312	2314	rBV2	42176	53858	2.98%	0.198%
86	15.212	2314	2317	2321	rVV2	77700	98223	5.44%	0.361%
87	15.280	2324	2328	2332	rVB4	32630	47155	2.61%	0.173%
88	15.377	2340	2344	2347	rBV6	14099	19872	1.10%	0.073%
89	15.493	2357	2363	2369	rBV	163663	242325	13.42%	0.891%
90	15.560	2370	2374	2380	rVV	70688	136965	7.59%	0.504%
91	15.615	2380	2383	2397	rVB7	32973	95870	5.31%	0.353%
92	15.889	2424	2428	2433	rVB4	13172	26163	1.45%	0.096%
93	16.090	2457	2461	2468	rBV8	18186	49599	2.75%	0.182%
94	16.377	2502	2508	2516	rVB	86331	142336	7.88%	0.523%

Sum of corrected areas: 27194424

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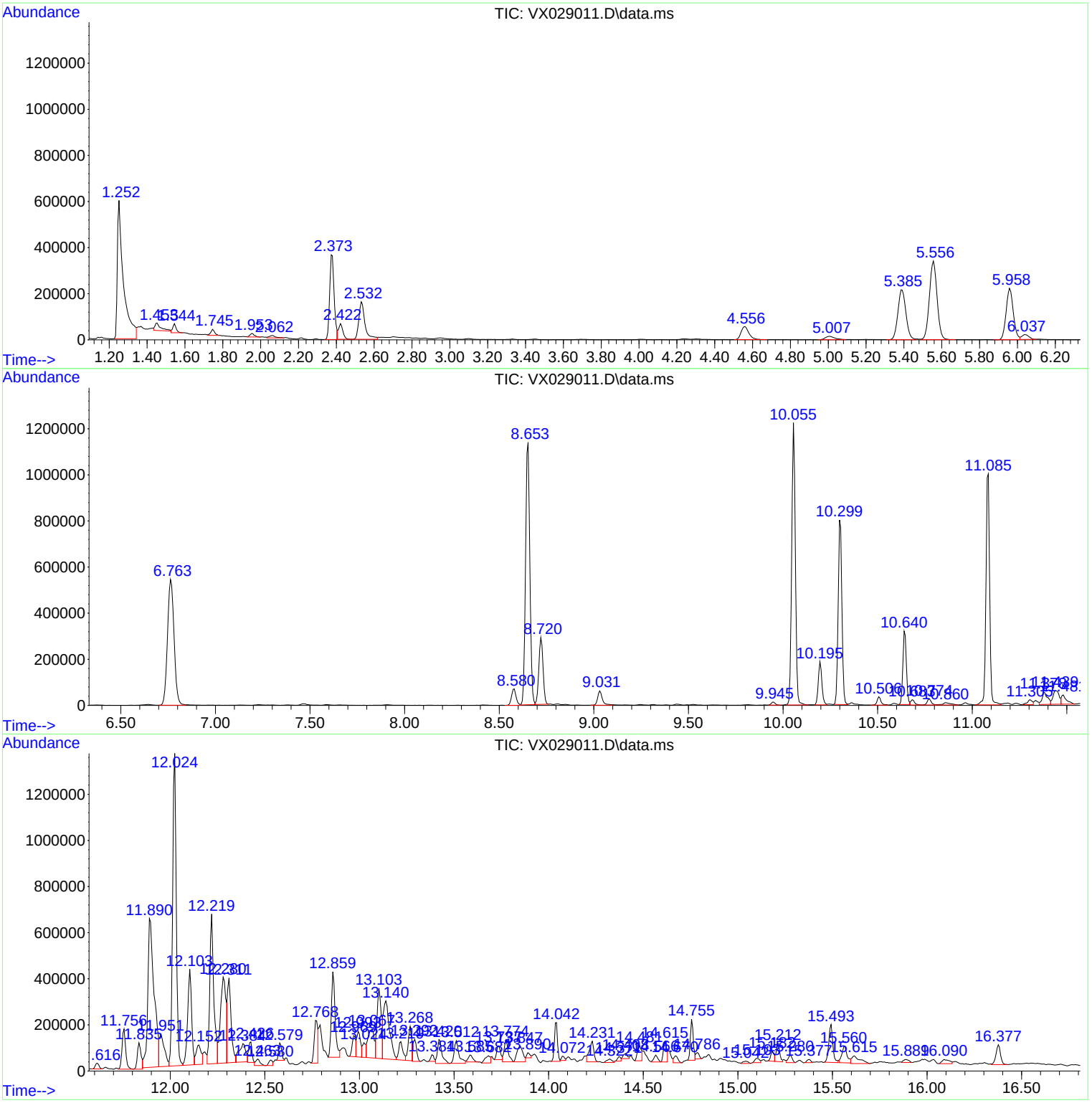
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
4922A

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\VX052622\
Data File : VX029011.D
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Operator : JC/MD
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Instrument :
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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

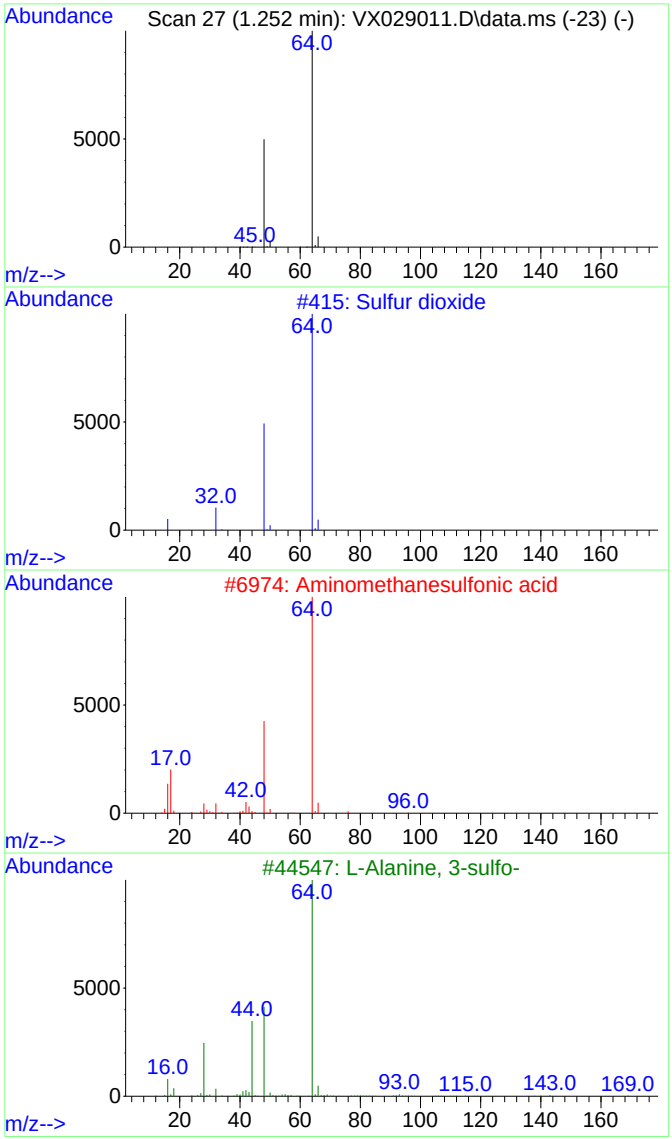
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

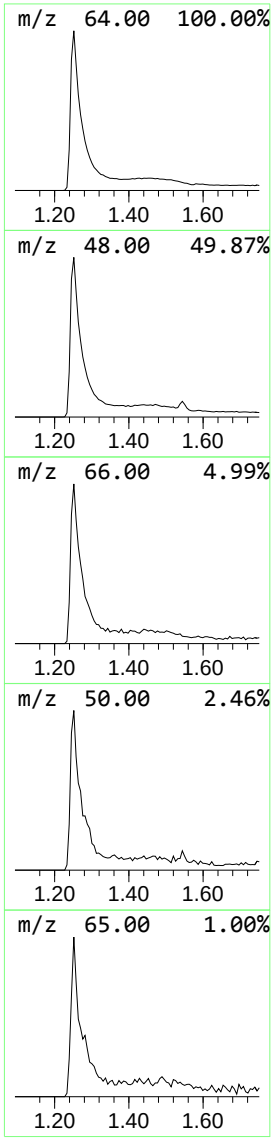
****

Peak Number 1 Sulfur dioxide Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.252	69.15 ug/l	1331100	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Sulfur dioxide	64	O2S	007446-09-5	83
2	Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	74
3	L-Alanine, 3-sulfo-	169	C3H7NO5S	000498-40-8	9
4	Ethene, 1,1-difluoro-	64	C2H2F2	000075-38-7	4
5	Ethyl Chloride	64	C2H5Cl	000075-00-3	3





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX052622\
Data File : VX029011.D
Acq On : 26 May 2022 19:55
Operator : JC/MD
Sample : N3059-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
4922A

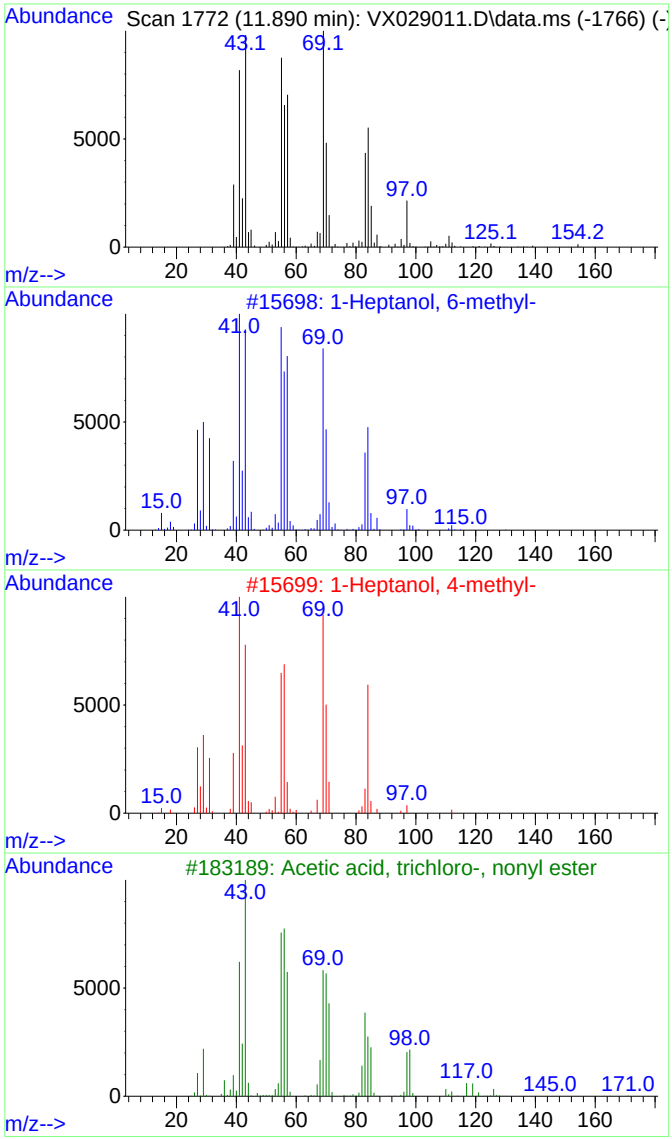
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

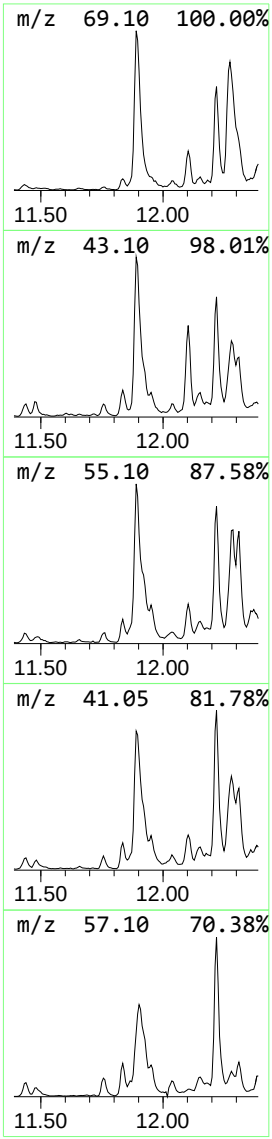
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 1-Heptanol, 6-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.890	41.36 ug/l	1493600	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heptanol, 6-methyl-	130	C8H18O	001653-40-3	86
2	1-Heptanol, 4-methyl-	130	C8H18O	000817-91-4	59
3	Acetic acid, trichloro-, nonyl e...	288	C11H19Cl3O2	065611-32-7	53
4	Cyclopropane, nonyl-	168	C12H24	074663-85-7	47
5	1-Pentene, 3-methyl-	84	C6H12	000760-20-3	47





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX052622\
Data File : VX029011.D
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Instrument :
MSVOA_X
ClientSampleId :
4922A

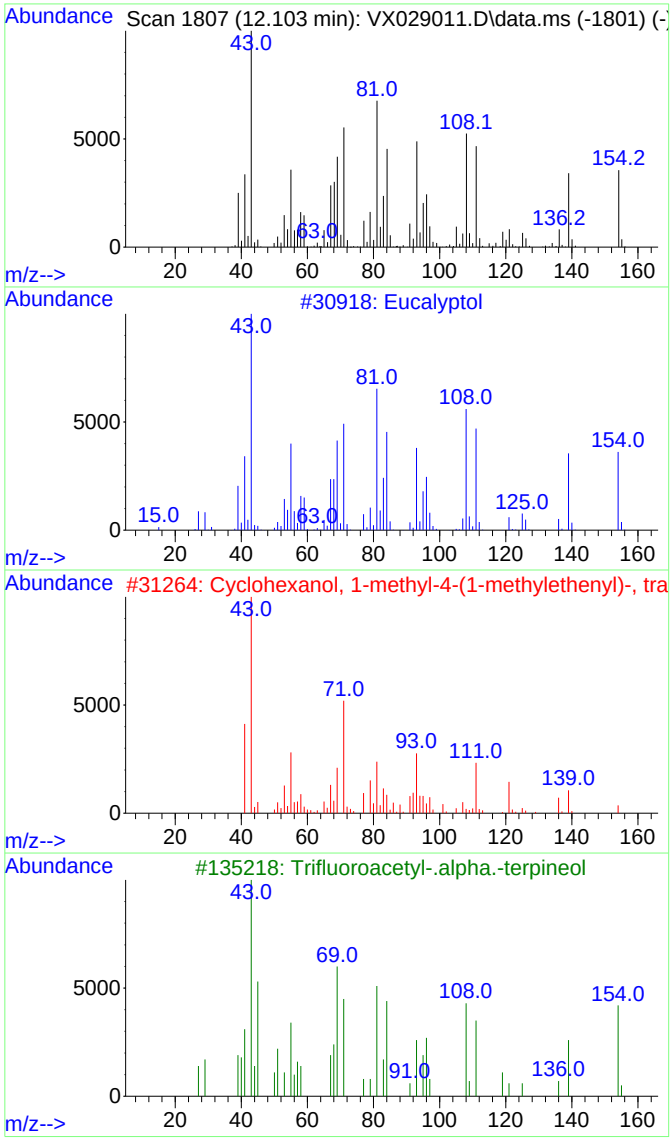
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

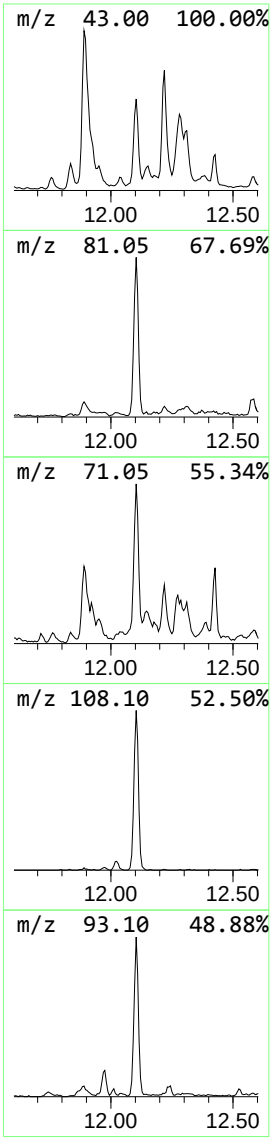
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Eucalyptol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.103	16.99 ug/l	613434	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eucalyptol	154	C10H18O	000470-82-6	99
2	Cyclohexanol, 1-methyl-4-(1-meth...	154	C10H18O	007299-40-3	47
3	Trifluoroacetyl-.alpha.-terpineol	250	C12H17F3O2	1000058-17-6	43
4	2-Cyclohexen-1-ol, 1-methyl-4-(1...	154	C10H18O	029803-82-5	35
5	5-Hepten-2-one, 6-methyl-	126	C8H14O	000110-93-0	22





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX052622\
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ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_X
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4922A

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

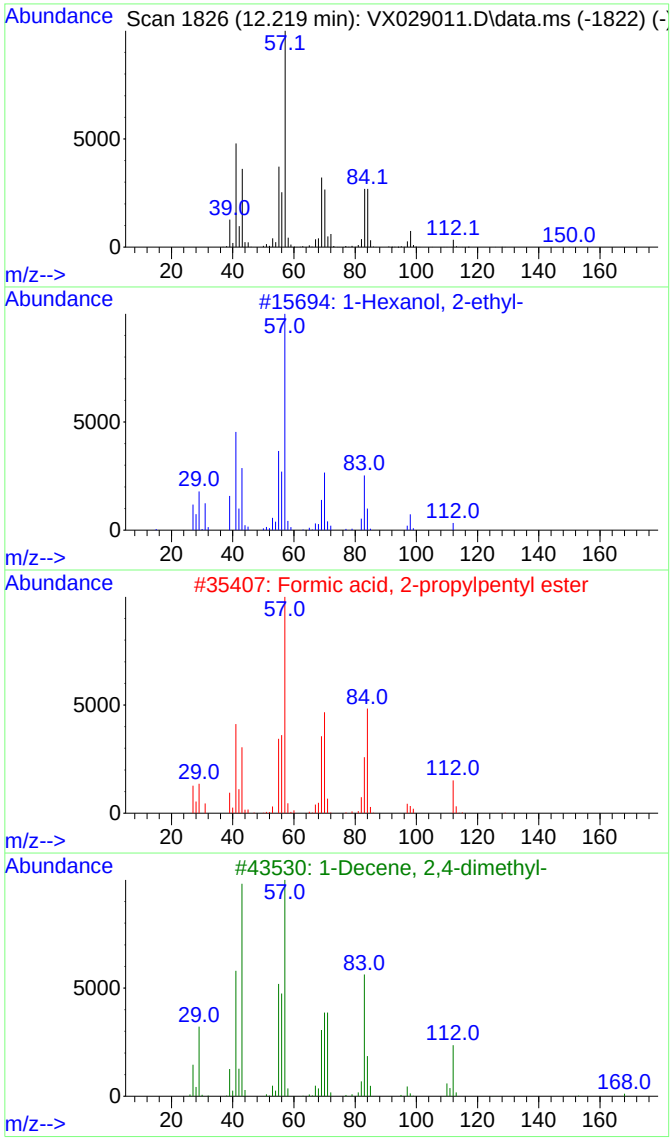
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

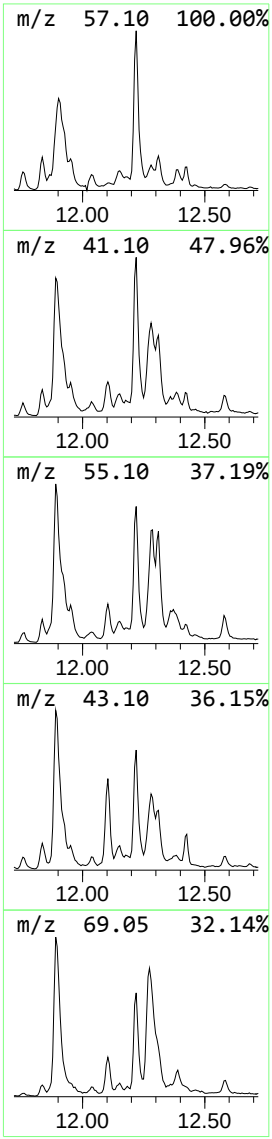
****

Peak Number 4 1-Hexanol, 2-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.219	26.47 ug/l	955869	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
2	Formic acid, 2-propylpentyl ester	158	C9H18O2	1000368-74-8	53
3	1-Decene, 2,4-dimethyl-	168	C12H24	055170-80-4	53
4	Propanoic acid, octyl ester	186	C11H22O2	000142-60-9	50
5	2-Propyl-1-Pentanol, trifluoroac...	226	C10H17F3O2	1000365-19-7	42





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX052622\
Data File : VX029011.D
Acq On : 26 May 2022 19:55
Operator : JC/MD
Sample : N3059-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
4922A

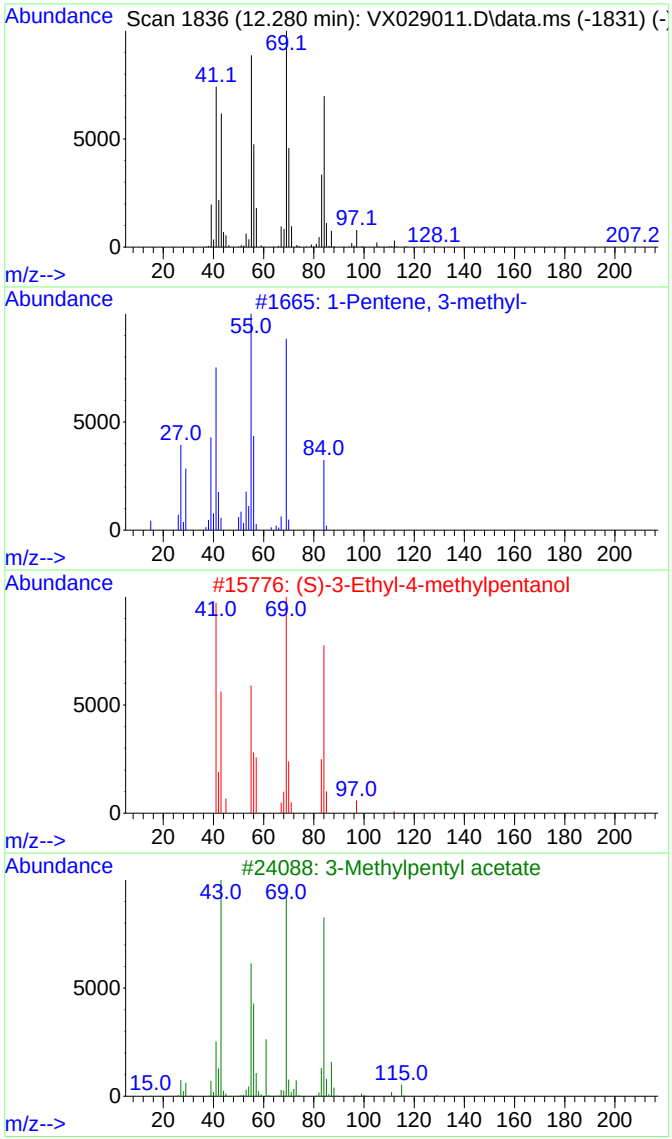
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

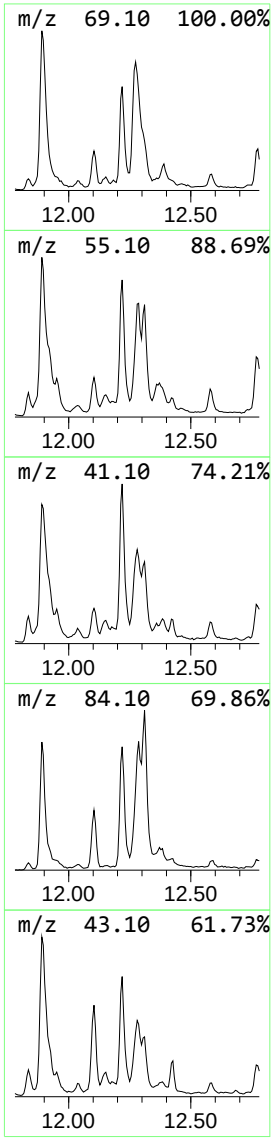
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 1-Pentene, 3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.280	21.21 ug/l	765775	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentene, 3-methyl-	84	C6H12	000760-20-3	53
2	(S)-3-Ethyl-4-methylpentanol	130	C8H18O	1000144-07-1	53
3	3-Methylpentyl acetate	144	C8H16O2	035897-13-3	53
4	Pentane, 3-methylene-	84	C6H12	000760-21-4	50
5	1,2-Cyclohexanedione	112	C6H8O2	000765-87-7	49





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX052622\
Data File : VX029011.D
Acq On : 26 May 2022 19:55
Operator : JC/MD
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ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
4922A

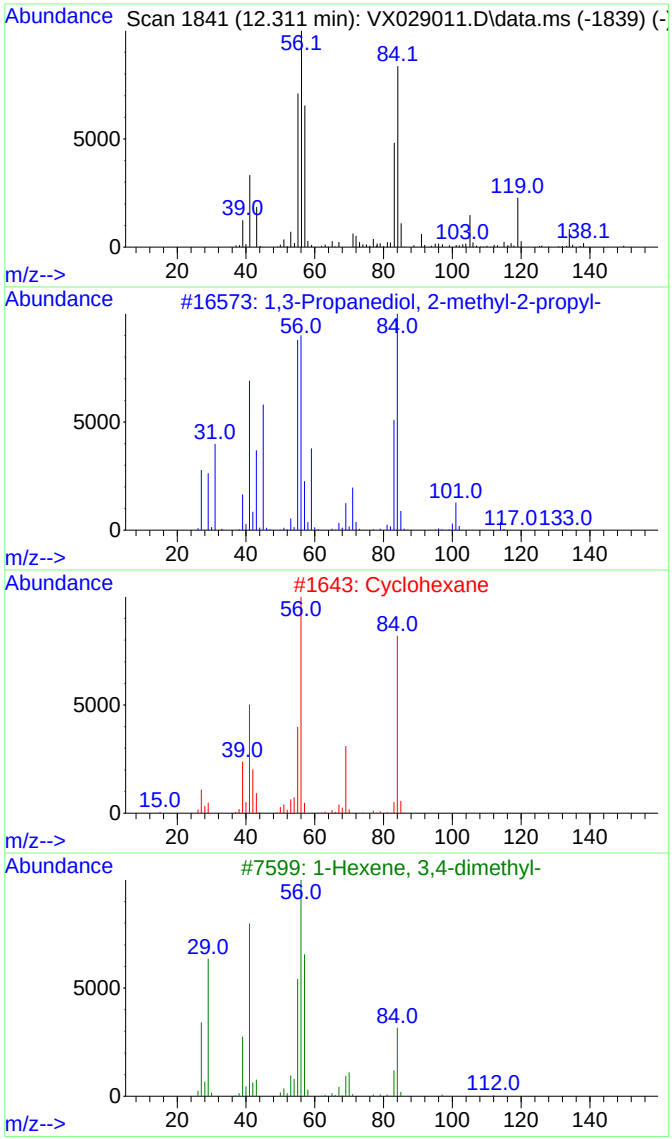
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

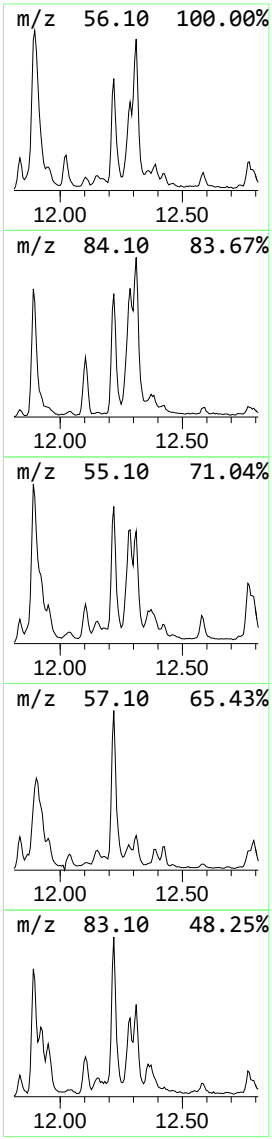
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 1,3-Propanediol, 2-methyl-2... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.311	13.58 ug/l	490311	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Propanediol, 2-methyl-2-propyl-	132	C7H16O2	000078-26-2	50
2	Cyclohexane	84	C6H12	000110-82-7	50
3	1-Hexene, 3,4-dimethyl-	112	C8H16	016745-94-1	45
4	1-(1,1-Dioxo-tetrahydro-1.lambda...	267	C9H17NO4S2	1000297-28-4	43
5	1-Hexanol, 2-(hydroxymethyl)-	132	C7H16O2	1000326-00-5	36





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX052622\
Data File : VX029011.D
Acq On : 26 May 2022 19:55
Operator : JC/MD
Sample : N3059-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
4922A

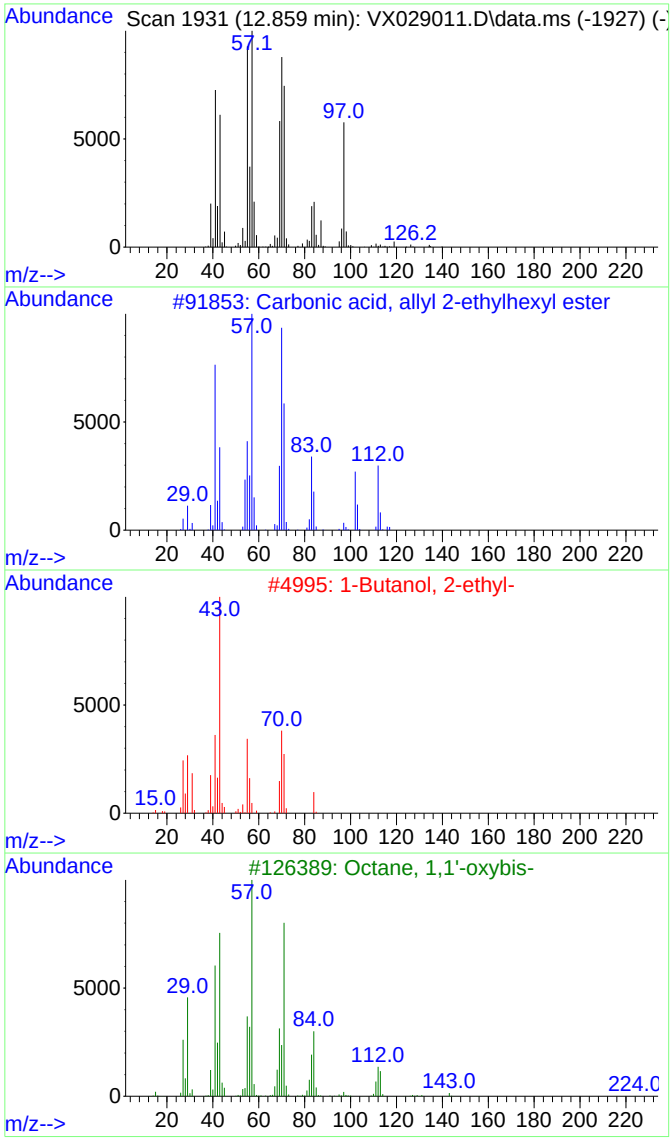
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

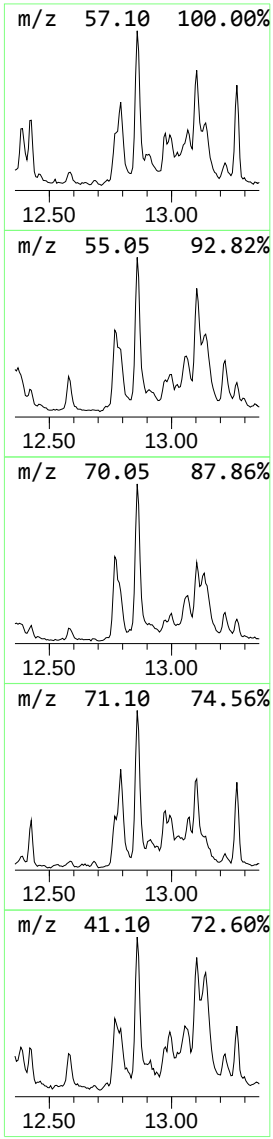
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Carbonic acid, allyl 2-ethy... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.859	14.39 ug/l	519774	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Carbonic acid, allyl 2-ethylhexy...	214	C12H22O3	1000357-80-6	64
2	1-Butanol, 2-ethyl-	102	C6H14O	000097-95-0	53
3	Octane, 1,1'-oxybis-	242	C16H34O	000629-82-3	35
4	Carbonic acid, octyl vinyl ester	200	C11H20O3	1000383-25-5	35
5	Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	27





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX052622\
Data File : VX029011.D
Acq On : 26 May 2022 19:55
Operator : JC/MD
Sample : N3059-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
4922A

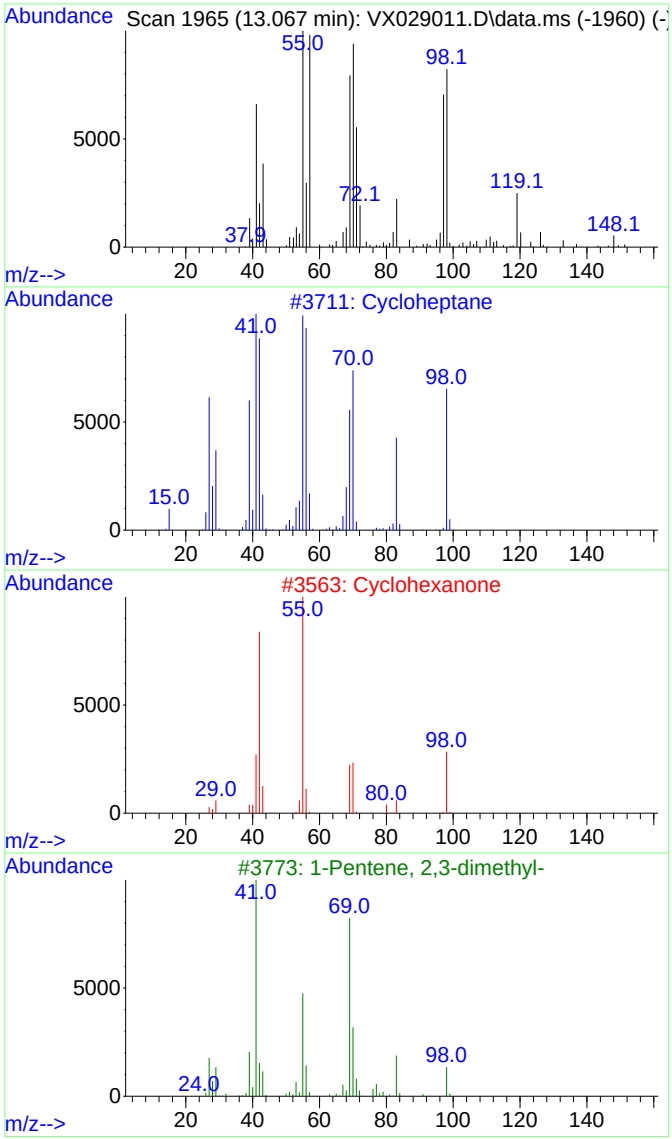
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

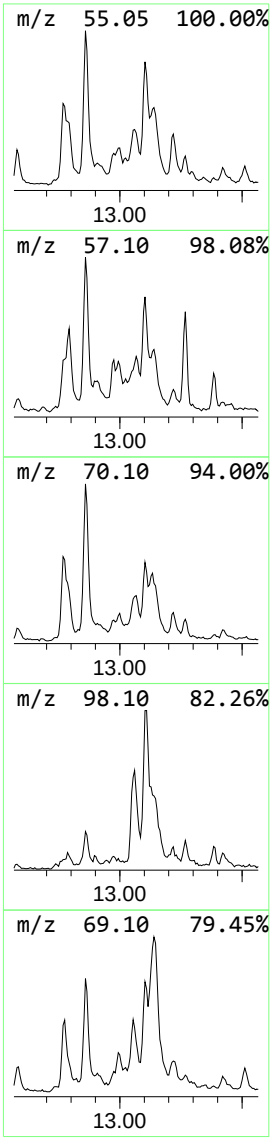
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 unknown13.067 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.067	8.01 ug/l	289281	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cycloheptane	98	C7H14	000291-64-5	47
2	Cyclohexanone	98	C6H10O	000108-94-1	43
3	1-Pentene, 2,3-dimethyl-	98	C7H14	003404-72-6	38
4	1-Hexanol, 4-methyl-	116	C7H16O	000818-49-5	35
5	Cyclohexane, 1,2,3-trimethyl-	126	C9H18	001678-97-3	35





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX052622\
Data File : VX029011.D
Acq On : 26 May 2022 19:55
Operator : JC/MD
Sample : N3059-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
4922A

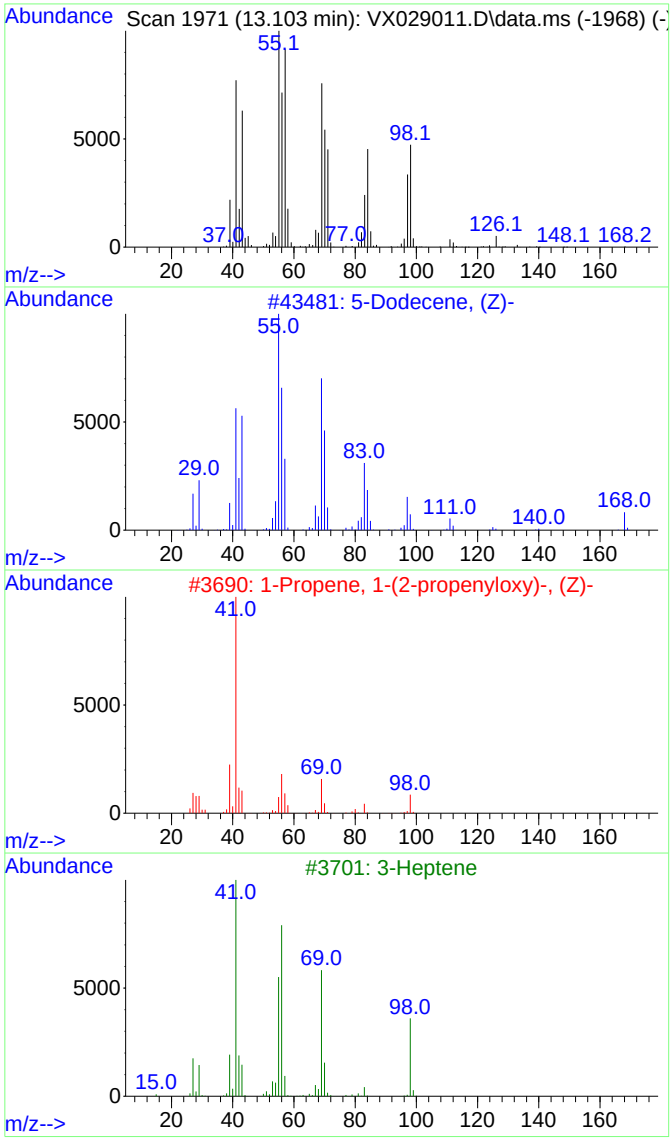
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

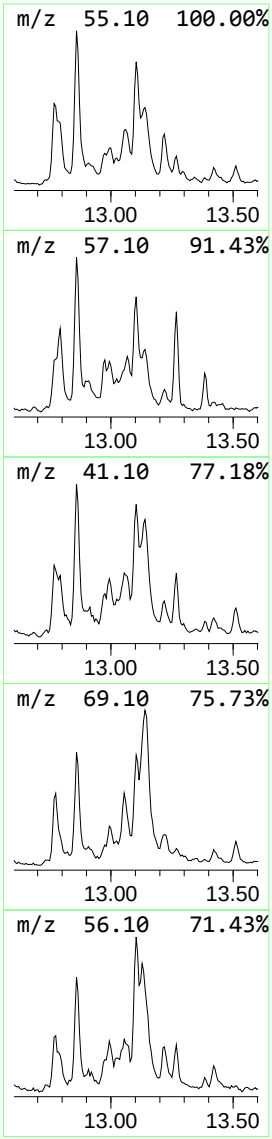
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 5-Dodecene, (Z)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.103	13.27 ug/l	479210	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Dodecene, (Z)-	168	C12H24	007206-28-2	50
2	1-Propene, 1-(2-propenyloxy)-, (Z)-	98	C6H10O	061142-12-9	43
3	3-Heptene	98	C7H14	000592-78-9	38
4	1-Octene, 3-methyl-	126	C9H18	013151-08-1	38
5	(Z)-2-Heptene	98	C7H14	006443-92-1	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX052622\
Data File : VX029011.D
Acq On : 26 May 2022 19:55
Operator : JC/MD
Sample : N3059-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
4922A

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

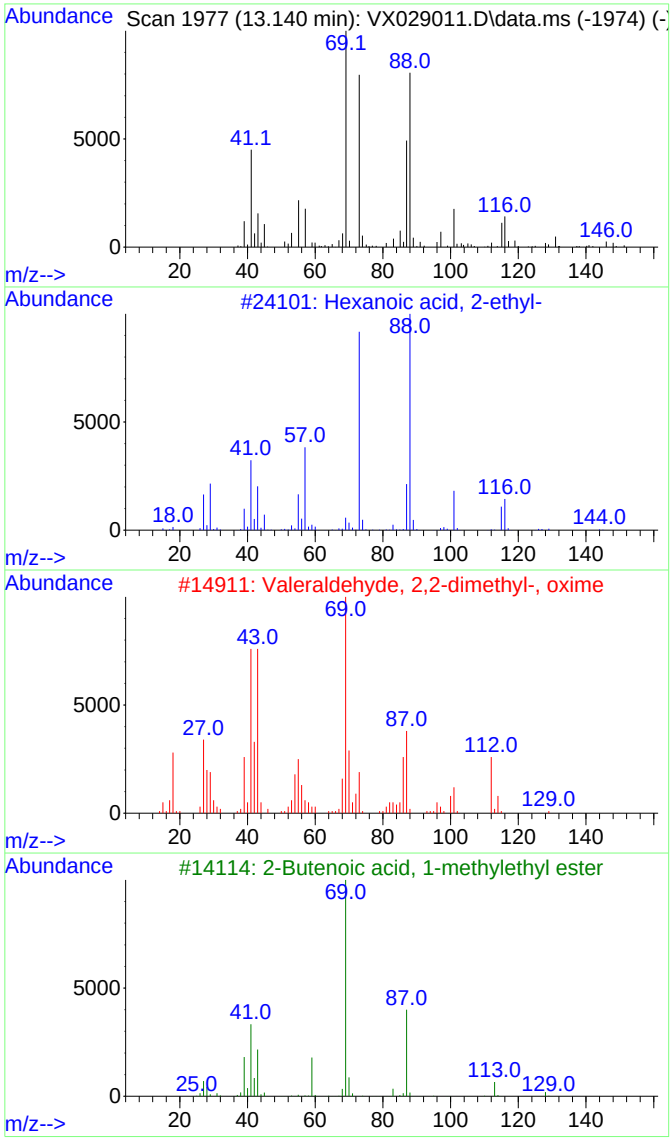
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

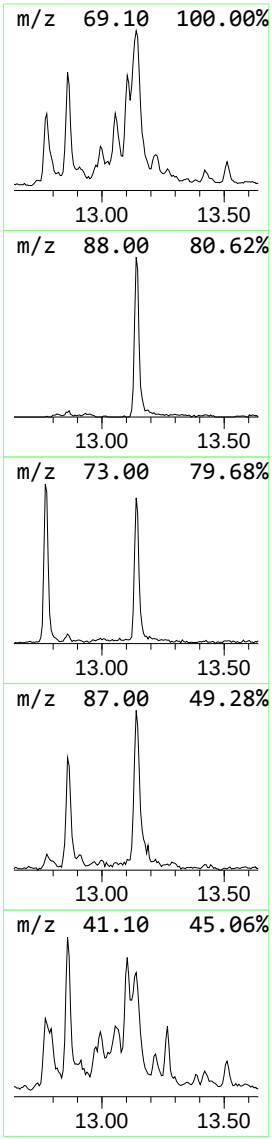
****

Peak Number 10 unknown13.140 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.140	15.31 ug/l	552669	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	43
2	Valeraldehyde, 2,2-dimethyl-, oxime	129	C7H15NO	016519-70-3	38
3	2-Butenoic acid, 1-methylethyl e...	128	C7H12O2	018060-77-0	38
4	Butanoic acid, 2-ethyl-, 1,2,3-p...	386	C21H38O6	056554-54-2	32
5	Acetic acid, diethyl-	116	C6H12O2	000088-09-5	32





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX052622\
Data File : VX029011.D
Acq On : 26 May 2022 19:55
Operator : JC/MD
Sample : N3059-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
4922A

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
Sulfur dioxide	1.252	69.2	ug/l	1331100	1	5.556	962527	50.0
1-Heptanol, 6-m...	11.890	41.4	ug/l	1493600	4	12.024	1805480	50.0
Eucalyptol	12.103	17.0	ug/l	613434	4	12.024	1805480	50.0
1-Hexanol, 2-et...	12.219	26.5	ug/l	955869	4	12.024	1805480	50.0
1-Pentene, 3-me...	12.280	21.2	ug/l	765775	4	12.024	1805480	50.0
1,3-Propanediol...	12.311	13.6	ug/l	490311	4	12.024	1805480	50.0
Carbonic acid, ...	12.859	14.4	ug/l	519774	4	12.024	1805480	50.0
unknown13.067	13.067	8.0	ug/l	289281	4	12.024	1805480	50.0
5-Dodecene, (Z)-	13.103	13.3	ug/l	479210	4	12.024	1805480	50.0
unknown13.140	13.140	15.3	ug/l	552669	4	12.024	1805480	50.0