LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022822\
Data File : VX027234.D
Acq On : 28 Feb 2022 20:44
Operator : JC/MD
Sample : N1688-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-22

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\82X022322W.M
Title : SW846 8260

Signal : TIC: VX027234.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	22	27	42	rBV3	26942	58950	13.47%	1.204%
2	1.374	44	47	53	rVB	8367	8174	1.87%	0.167%
3	1.752	104	109	118	rVB	40300	55235	12.62%	1.128%
4	1.959	138	143	152	rVB2	8121	12400	2.83%	0.253%
5	2.221	181	186	193	rVB	3741	5935	1.36%	0.121%
6	2.343	199	206	210	rBV2	2575	5013	1.15%	0.102%
7	2.782	269	278	282	rBV2	9318	21242	4.86%	0.434%
8	2.831	282	286	288	rVV2	9152	16559	3.78%	0.338%
9	2.873	288	293	301	rVV	20226	43102	9.85%	0.880%
10	2.959	301	307	323	rVV	26464	58096	13.28%	1.187%
11	3.117	323	333	344	rVB3	34300	85738	19.60%	1.751%
12	3.763	428	439	446	rBV6	5280	17336	3.96%	0.354%
13	3.843	446	452	460	rVB3	2379	5343	1.22%	0.109%
14	4.312	518	529	541	rVB	44445	125646	28.72%	2.566%
15	4.434	541	549	559	rVB4	2282	6726	1.54%	0.137%
16	5.190	668	673	684	rVB5	2057	5980	1.37%	0.122%
17	5.391	695	706	714	rBV	50343	148925	34.04%	3.042%
18	5.477	714	720	726	rVV	37101	110202	25.19%	2.251%
19	5.556	726	733	757	rVB	86632	260243	59.48%	5.315%
20	5.848	773	781	790	rVB6	1989	6221	1.42%	0.127%
21	5.964	790	800	806	rBV	55304	141319	32.30%	2.886%
22	6.043	806	813	826	rVV	89589	236550	54.07%	4.831%
23	6.251	826	847	856	rVV9	6207	34752	7.94%	0.710%
24	6.361	856	865	875	rVB7	4023	12733	2.91%	0.260%
25	6.769	921	932	943	rBV	134088	324659	74.21%	6.631%
26	6.848	943	945	954	rVB4	3113	6833	1.56%	0.140%
27	7.171	992	998	1006	rBV3	4635	9578	2.19%	0.196%
28	7.379	1022	1032	1040	rBV3	14578	38741	8.85%	0.791%
29	7.885	1103	1115	1122	rBV6	4544	13882	3.17%	0.284%
30	7.988	1122	1132	1139	rVB2	2603	5739	1.31%	0.117%
31	8.110	1145	1152	1154	rBV2	5048	9208	2.10%	0.188%
32	8.147	1154	1158	1164	rVB	8745	15158	3.46%	0.310%
33	8.427	1196	1204	1211	rBV2	4735	8360	1.91%	0.171%
34	8.506	1211	1217	1228	rVB2	11855	19848	4.54%	0.405%
35	8.653	1228	1241	1249	rBV	280970	437516	100.00%	8.936%
36	8.958	1282	1291	1299	rVB3	13298	24941	5.70%	0.509%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022822\
Data File : VX027234.D
Acq On : 28 Feb 2022 20:44
Operator : JC/MD
Sample : N1688-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-22

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\82X022322W.M
Title : SW846 8260

37	9.049	1299	1306	1312	rBV	12825	20811	4.76%	0.425%
38	9.165	1319	1325	1332	rBV	6304	9792	2.24%	0.200%
39	9.458	1364	1373	1378	rBV3	9907	19295	4.41%	0.394%
40	9.945	1449	1453	1461	rVB2	3390	6073	1.39%	0.124%
41	10.055	1465	1471	1477	rBV	286394	382032	87.32%	7.803%
42	10.195	1490	1494	1499	rVB	28949	36635	8.37%	0.748%
43	10.305	1509	1512	1518	rVB2	3835	5129	1.17%	0.105%
44	10.963	1615	1620	1626	rVB	53148	65145	14.89%	1.331%
45	11.079	1634	1639	1650	rVB	229552	296069	67.67%	6.047%
46	11.305	1671	1676	1685	rVB	42703	51776	11.83%	1.057%
47	11.616	1721	1727	1734	rVB	11979	15125	3.46%	0.309%
48	11.756	1746	1750	1755	rVB	5852	7522	1.72%	0.154%
49	11.866	1764	1768	1770	rBV	8082	9784	2.24%	0.200%
50	11.890	1770	1772	1777	rVB	8283	9243	2.11%	0.189%
51	12.024	1788	1794	1801	rBV	300771	372103	85.05%	7.600%
52	12.091	1801	1805	1814	rVB	5532	7350	1.68%	0.150%
53	12.177	1816	1819	1825	rBV	4124	4874	1.11%	0.100%
54	12.244	1825	1830	1837	rBV	243285	294524	67.32%	6.015%
55	12.317	1837	1842	1848	rVB2	9529	13967	3.19%	0.285%
56	12.408	1852	1857	1862	rBV3	12895	16753	3.83%	0.342%
57	12.463	1862	1866	1871	rVB	31116	35788	8.18%	0.731%
58	12.530	1873	1877	1884	rBV	4163	5554	1.27%	0.113%
59	12.615	1884	1891	1893	rBV2	5136	7082	1.62%	0.145%
60	12.646	1893	1896	1900	rVV	27800	33706	7.70%	0.688%
61	12.701	1900	1905	1912	rVB	128061	154524	35.32%	3.156%
62	12.847	1924	1929	1933	rVB2	13268	17865	4.08%	0.365%
63	12.920	1933	1941	1945	rVV	60247	78995	18.06%	1.613%
64	12.963	1945	1948	1953	rVV	53144	64042	14.64%	1.308%
65	13.036	1955	1960	1966	rVB2	5179	7621	1.74%	0.156%
66	13.140	1972	1977	1979	rBV2	5873	7334	1.68%	0.150%
67	13.170	1979	1982	1991	rVB	21926	29953	6.85%	0.612%
68	13.250	1991	1995	1998	rBV2	3989	5934	1.36%	0.121%
69	13.292	1998	2002	2007	rVV	76173	95593	21.85%	1.952%
70	13.347	2007	2011	2013	rVV3	5876	9130	2.09%	0.186%
71	13.371	2013	2015	2019	rVV	4057	4678	1.07%	0.096%
72	13.420	2019	2023	2030	rVB	31841	41502	9.49%	0.848%
73	13.506	2030	2037	2040	rBV3	4117	6707	1.53%	0.137%
74	13.542	2040	2043	2046	rVV	12978	17567	4.02%	0.359%
75	13.579	2046	2049	2055	rVV3	17848	33350	7.62%	0.681%
76	13.640	2056	2059	2061	rVV2	10227	15511	3.55%	0.317%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022822\
Data File : VX027234.D
Acq On : 28 Feb 2022 20:44
Operator : JC/MD
Sample : N1688-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-22

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\82X022322W.M
Title : SW846 8260

77	13.664	2061	2063	2070	rVB2	14961	20185	4.61%	0.412%
78	13.774	2074	2081	2091	rBV	40403	55311	12.64%	1.130%
79	13.859	2091	2095	2100	rBV3	5874	9869	2.26%	0.202%
80	13.926	2100	2106	2110	rBV2	8944	11981	2.74%	0.245%
81	13.969	2110	2113	2118	rVB	4097	4867	1.11%	0.099%
82	14.079	2126	2131	2134	rBV	6211	7675	1.75%	0.157%
83	14.213	2148	2153	2163	rBV2	10442	16514	3.77%	0.337%
84	14.322	2167	2171	2175	rBV	5464	6494	1.48%	0.133%
85	14.371	2175	2179	2184	rVB	6462	8501	1.94%	0.174%
86	14.481	2194	2197	2202	rBV2	4775	6397	1.46%	0.131%
87	14.780	2242	2246	2254	rBV	26744	34998	8.00%	0.715%

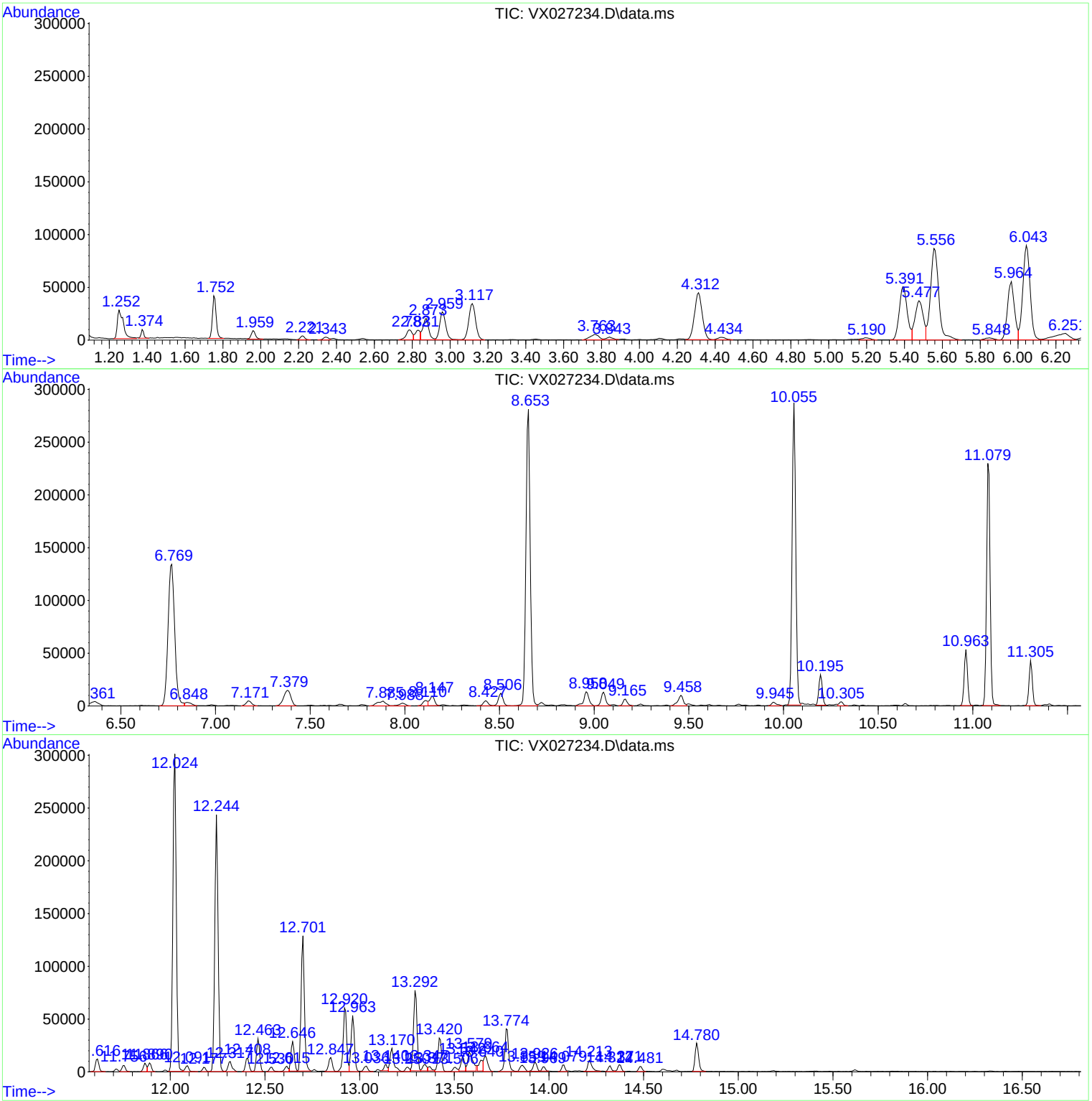
Sum of corrected areas: 4896113

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022822\
Data File : VX027234.D
Acq On : 28 Feb 2022 20:44
Operator : JC/MD
Sample : N1688-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-22

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022322W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022822\
Data File : VX027234.D
Acq On : 28 Feb 2022 20:44
Operator : JC/MD
Sample : N1688-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-22

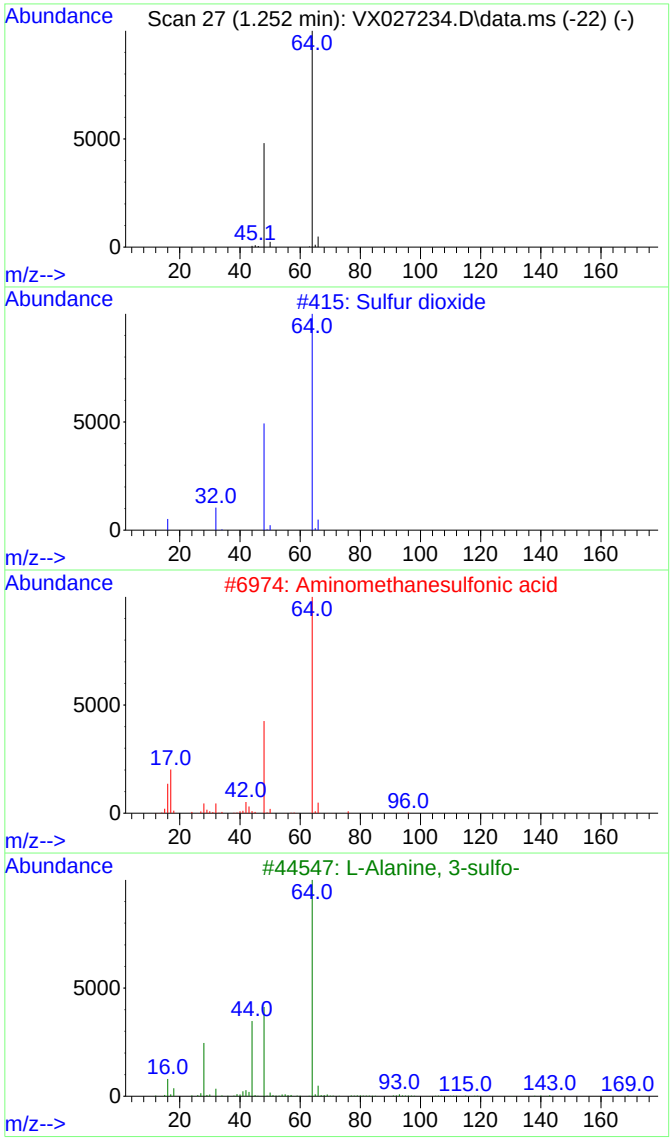
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022322W.M
Quant Title : SW846 8260

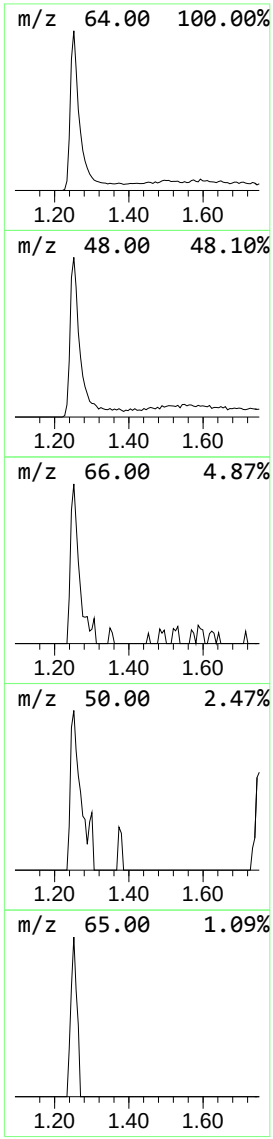
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Sulfur dioxide Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.252	11.33 ug/l	58950	Pentafluorobenzene	5.562

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Sulfur dioxide	64	O2S	007446-09-5	90
2	Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	64
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\VX022822\
Data File : VX027234.D
Acq On : 28 Feb 2022 20:44
Operator : JC/MD
Sample : N1688-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-22

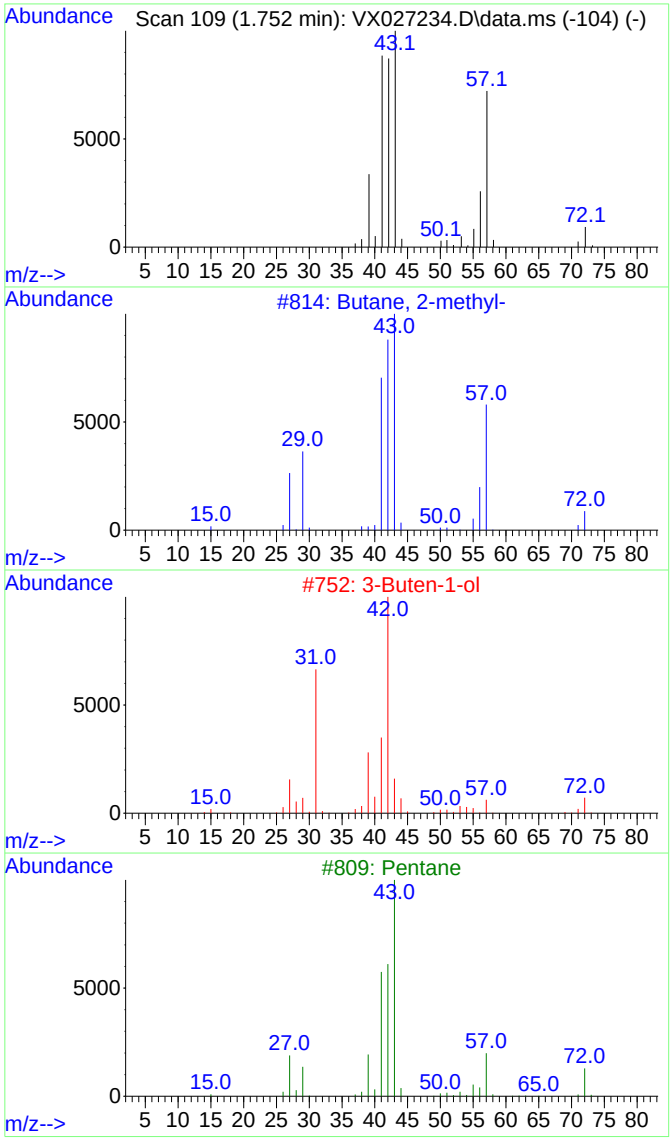
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022322W.M
Quant Title : SW846 8260

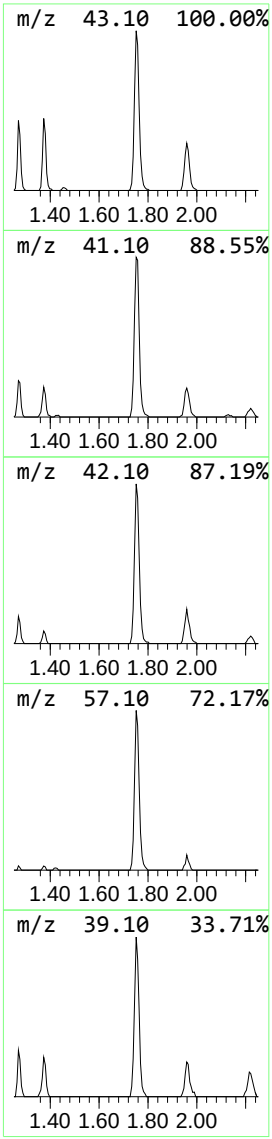
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.752	10.61 ug/l	55235	Pentafluorobenzene	5.562

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	91
2	3-Buten-1-ol	72	C4H8O	000627-27-0	47
3	Pentane	72	C5H12	000109-66-0	38
4	2(3H)-Furanone, dihydro-4-methyl-	100	C5H8O2	001679-49-8	33
5	1-Butene	56	C4H8	000106-98-9	14





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022822\
Data File : VX027234.D
Acq On : 28 Feb 2022 20:44
Operator : JC/MD
Sample : N1688-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-22

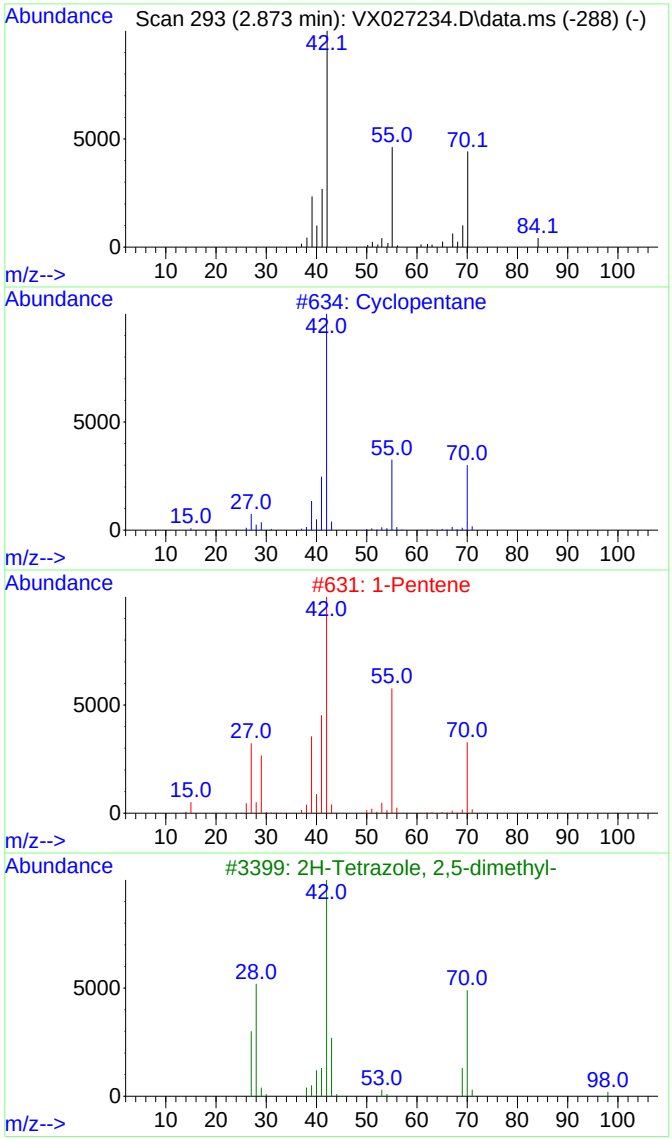
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022322W.M
Quant Title : SW846 8260

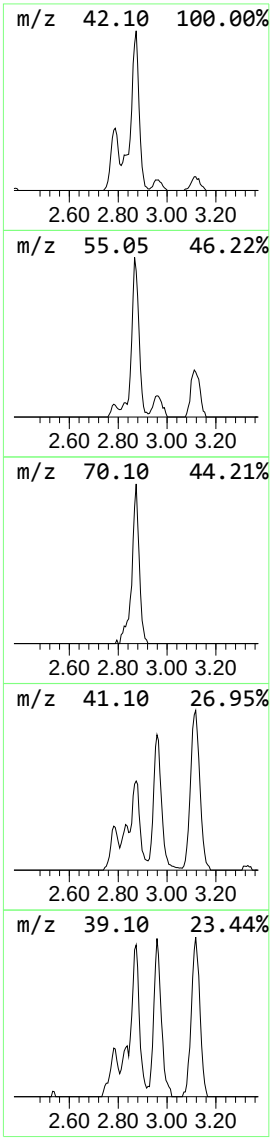
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Cyclopentane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.873	8.28 ug/l	43102	Pentafluorobenzene	5.562

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane	70	C5H10	000287-92-3	56
2	1-Pentene	70	C5H10	000109-67-1	45
3	2H-Tetrazole, 2,5-dimethyl-	98	C3H6N4	004135-93-7	40
4	3-Buten-1-ol	72	C4H8O	000627-27-0	25
5	2-Oxetanone, 4-methylene-	84	C4H4O2	000674-82-8	9





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022822\
Data File : VX027234.D
Acq On : 28 Feb 2022 20:44
Operator : JC/MD
Sample : N1688-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-22

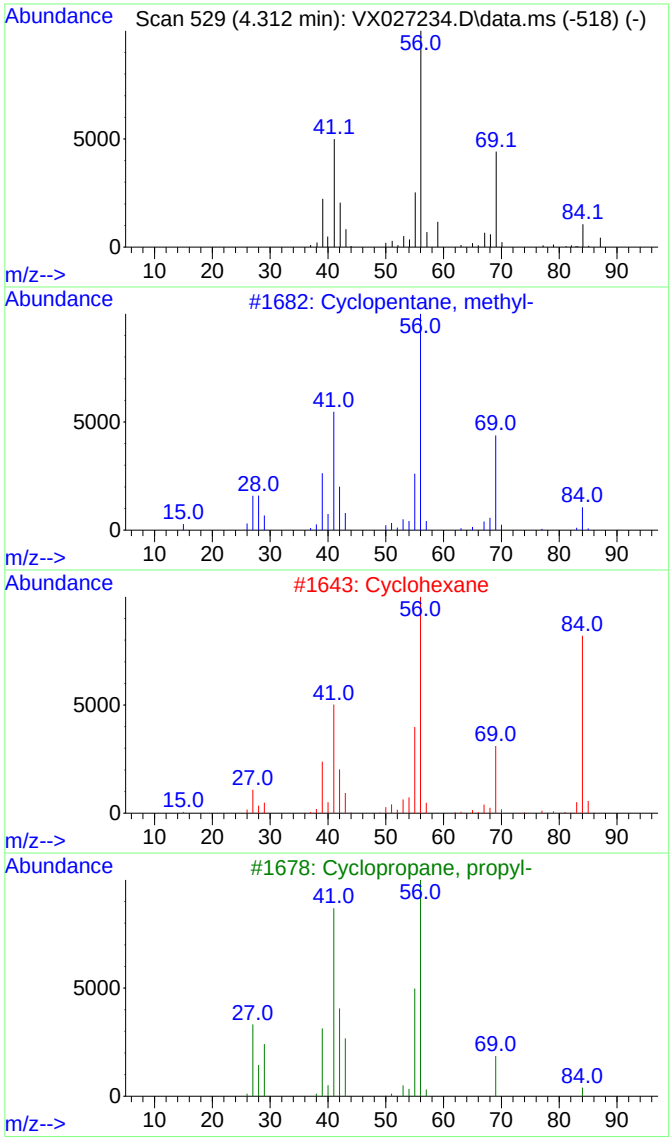
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022322W.M
Quant Title : SW846 8260

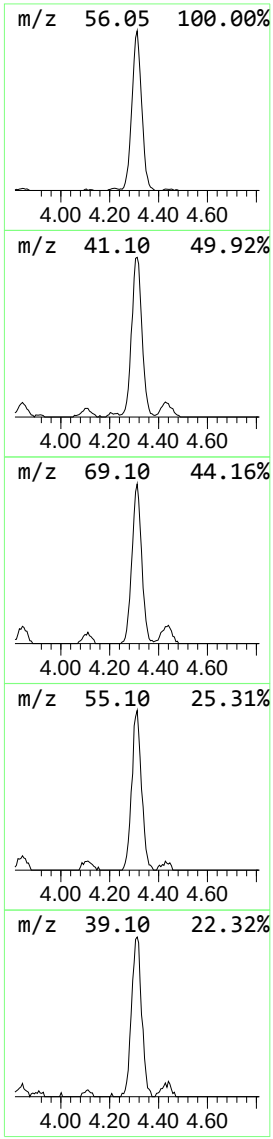
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Cyclopentane, methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.312	24.14 ug/l	125646	Pentafluorobenzene	5.562

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	93
2	Cyclohexane	84	C6H12	000110-82-7	64
3	Cyclopropane, propyl-	84	C6H12	002415-72-7	59
4	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	59
5	Cyclobutane, ethyl-	84	C6H12	004806-61-5	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022822\
Data File : VX027234.D
Acq On : 28 Feb 2022 20:44
Operator : JC/MD
Sample : N1688-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-22

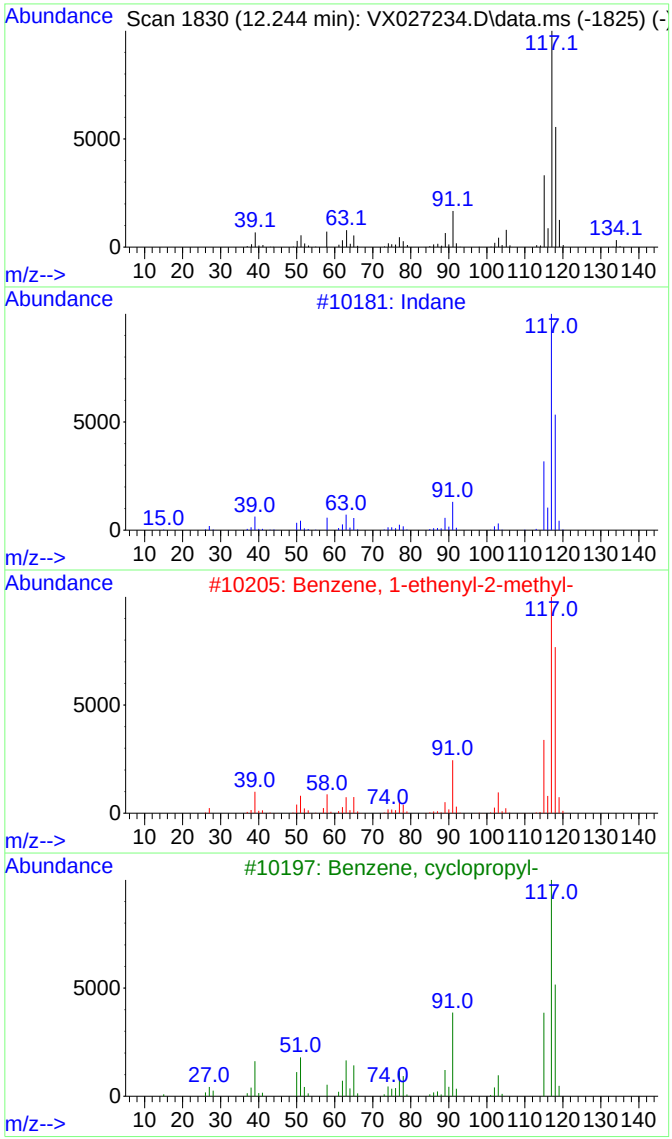
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022322W.M
Quant Title : SW846 8260

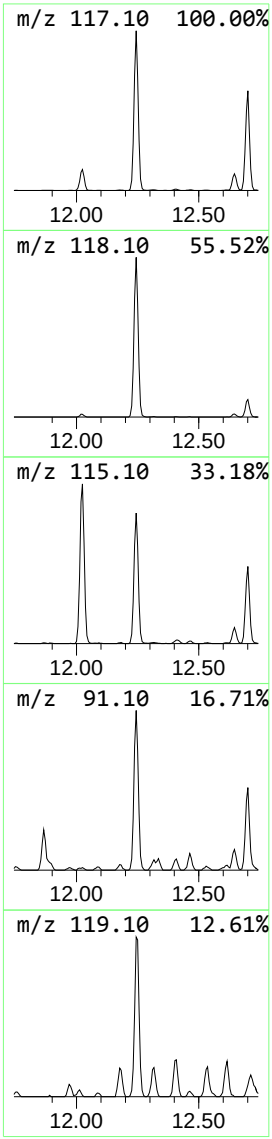
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	39.58 ug/l	294524	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	81
2	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	81
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	81
4	Benzene, 2-propenyl-	118	C9H10	000300-57-2	74
5	Deltacyclene	118	C9H10	007785-10-6	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022822\
Data File : VX027234.D
Acq On : 28 Feb 2022 20:44
Operator : JC/MD
Sample : N1688-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-22

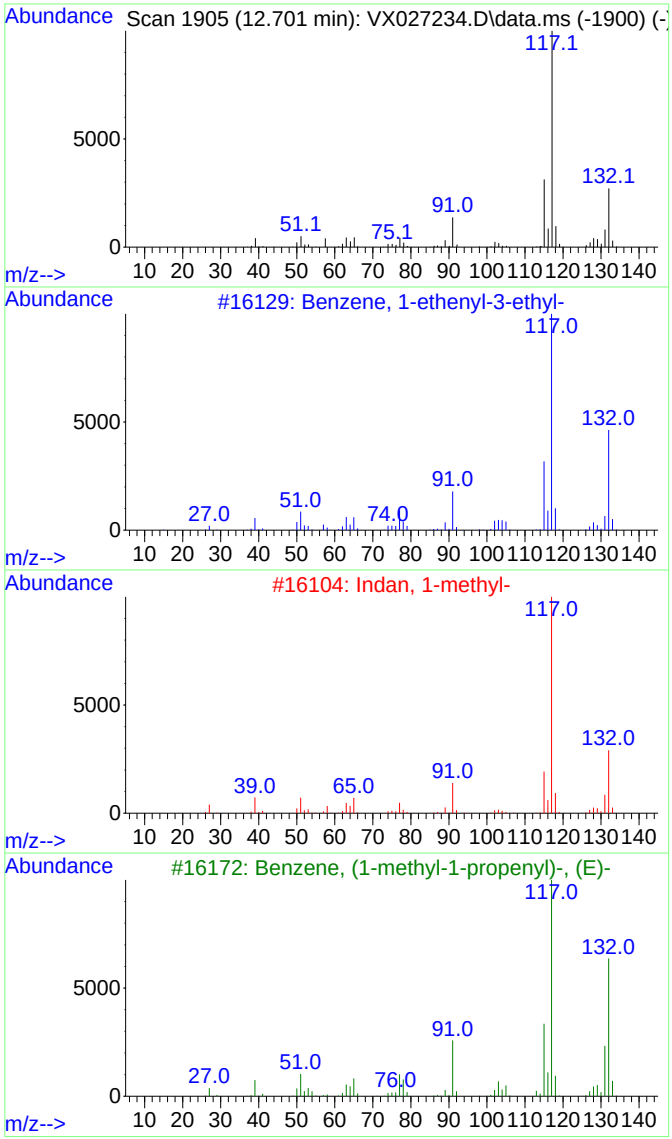
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022322W.M
Quant Title : SW846 8260

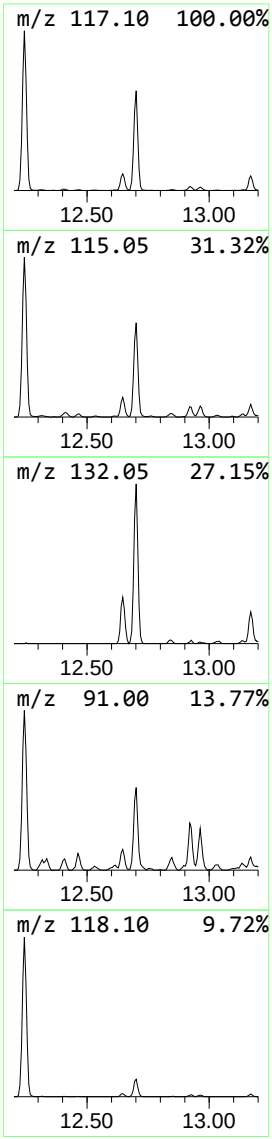
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	20.76 ug/l	154524	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	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	90
4	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
5	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022822\
Data File : VX027234.D
Acq On : 28 Feb 2022 20:44
Operator : JC/MD
Sample : N1688-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-22

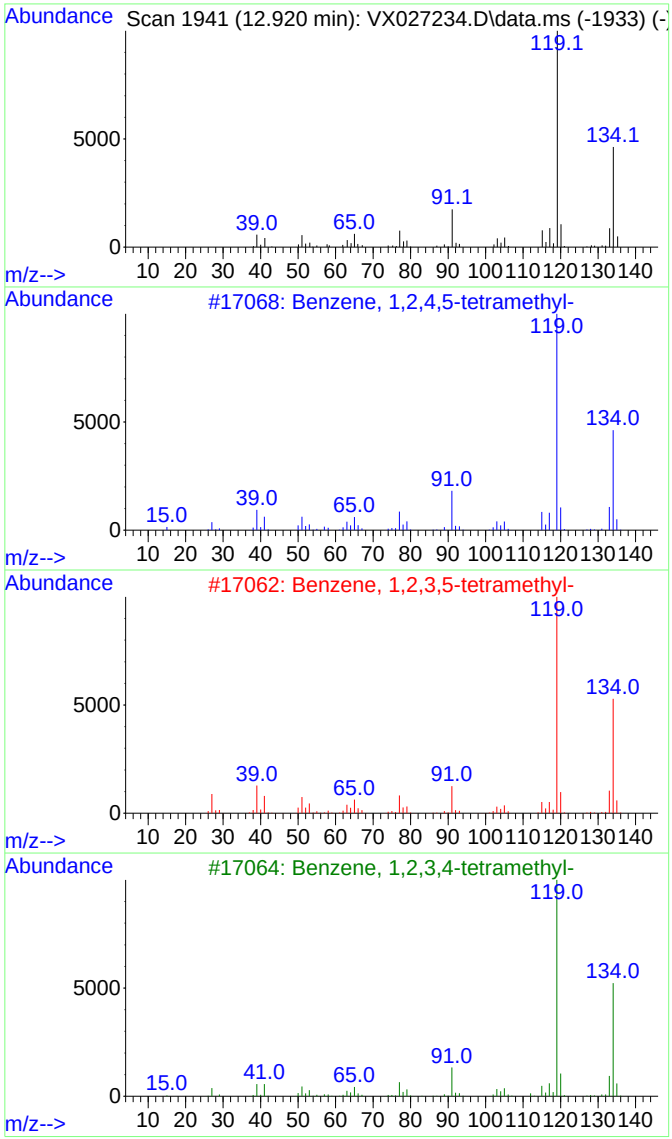
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022322W.M
Quant Title : SW846 8260

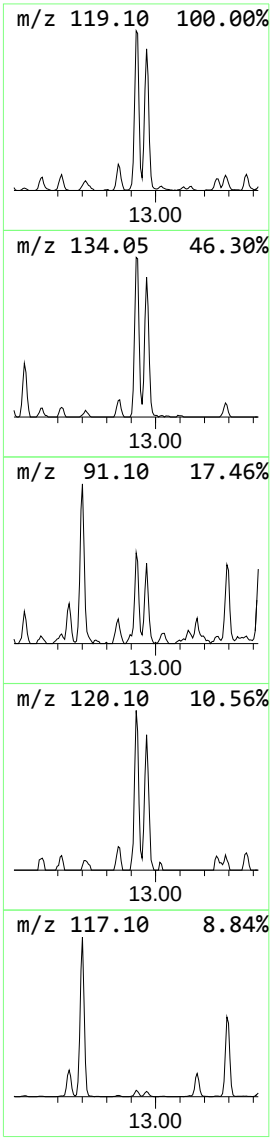
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.920	10.61 ug/l	78995	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
5	o-Cymene	134	C10H14	000527-84-4	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022822\
Data File : VX027234.D
Acq On : 28 Feb 2022 20:44
Operator : JC/MD
Sample : N1688-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-22

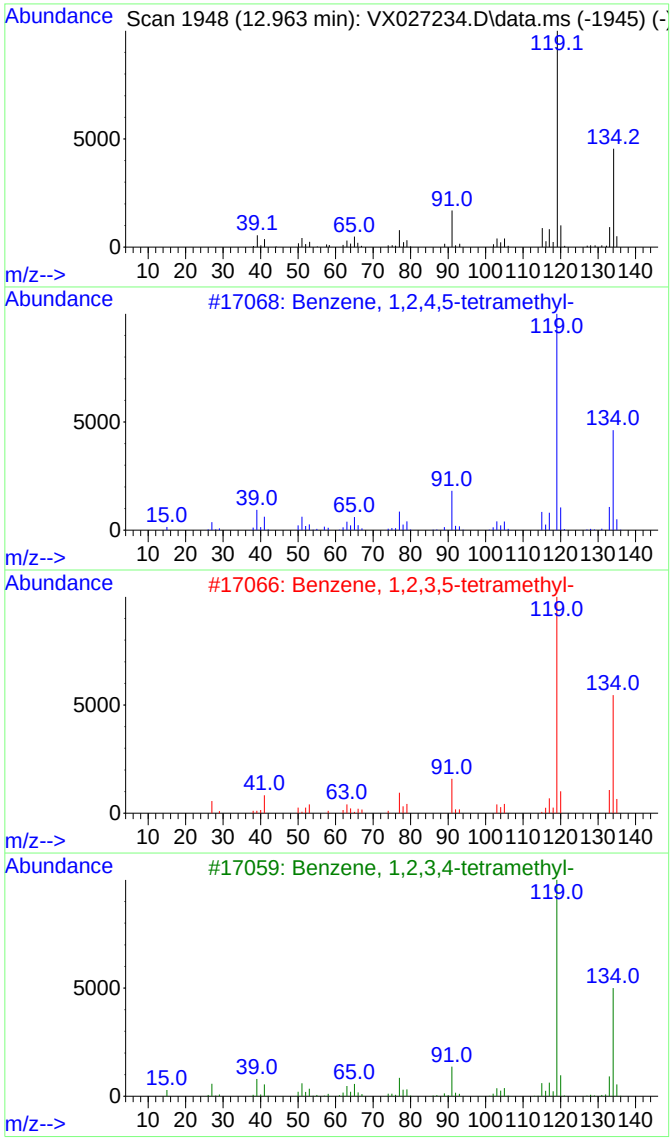
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022322W.M
Quant Title : SW846 8260

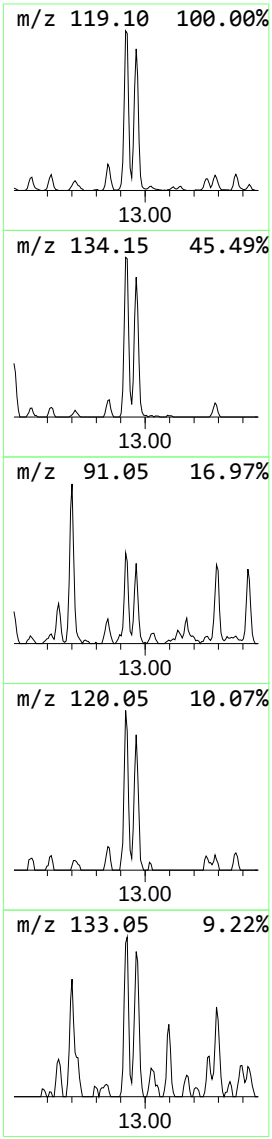
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.963	8.61 ug/l	64042	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022822\
Data File : VX027234.D
Acq On : 28 Feb 2022 20:44
Operator : JC/MD
Sample : N1688-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-22

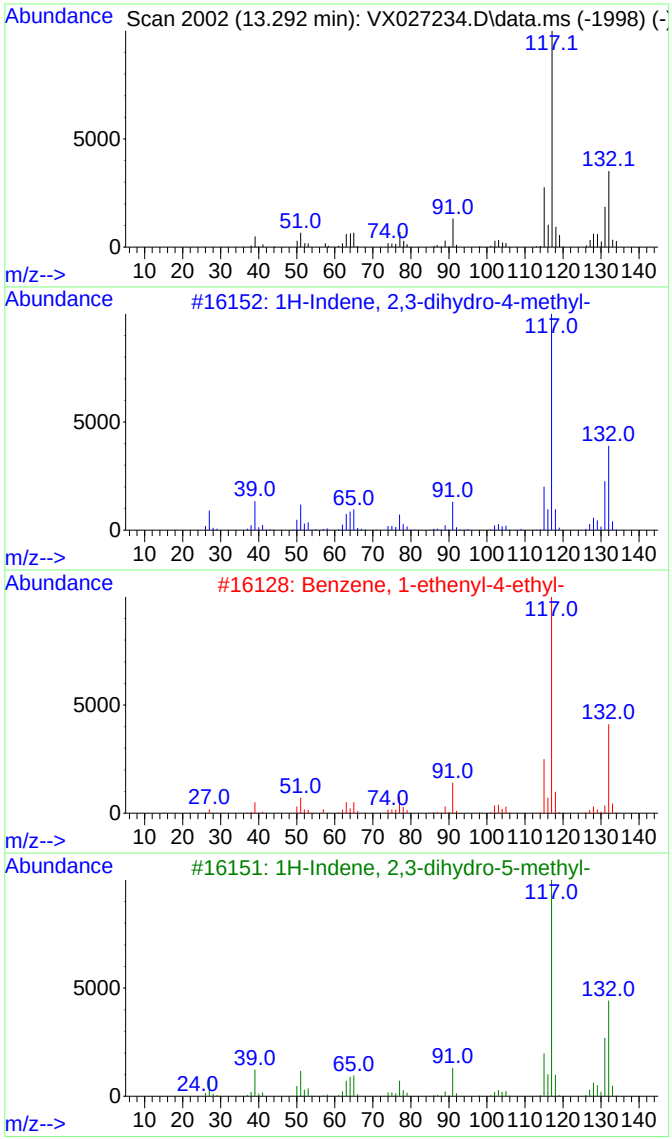
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022322W.M
Quant Title : SW846 8260

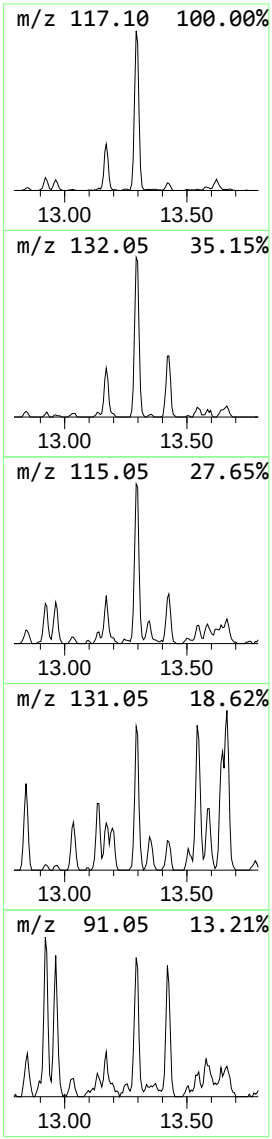
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	12.85 ug/l	95593	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	95
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022822\
Data File : VX027234.D
Acq On : 28 Feb 2022 20:44
Operator : JC/MD
Sample : N1688-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-22

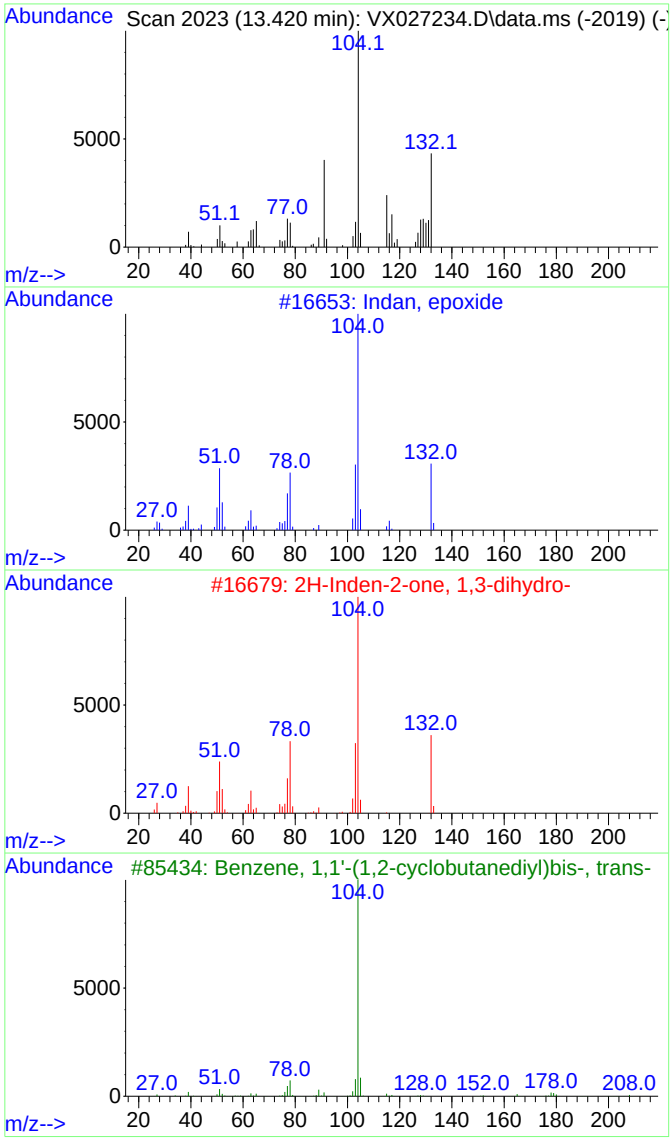
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022322W.M
Quant Title : SW846 8260

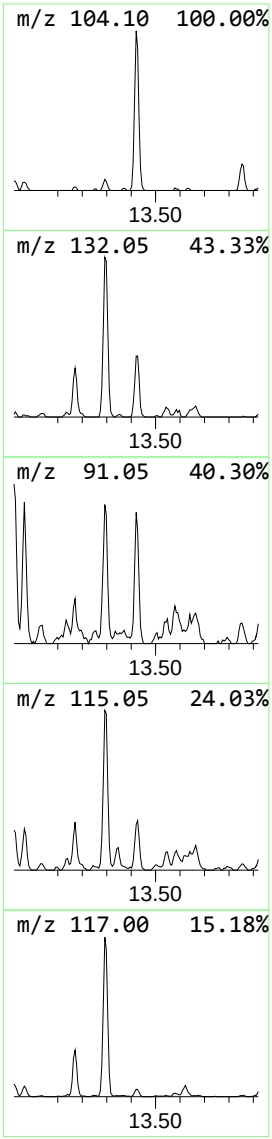
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 unknown13.420 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.420	5.58 ug/l	41502	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, epoxide	132	C9H8O	000768-22-9	46
2	2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	43
3	Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	020071-09-4	27
4	Benzenemethanol, 2-methyl-	122	C8H10O	000089-95-2	27
5	Pentafluoropropionic acid, 2-phe...	268	C11H9F5O2	081089-48-7	27





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022822\
Data File : VX027234.D
Acq On : 28 Feb 2022 20:44
Operator : JC/MD
Sample : N1688-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-22

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022322W.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	11.3	ug/l	58950	1	5.562	260243	50.0
Butane, 2-methyl-	1.752	10.6	ug/l	55235	1	5.562	260243	50.0
Cyclopentane	2.873	8.3	ug/l	43102	1	5.562	260243	50.0
Cyclopentane, m...	4.312	24.1	ug/l	125646	1	5.562	260243	50.0
Indane	12.244	39.6	ug/l	294524	4	12.024	372103	50.0
Benzene, 1-ethe...	12.701	20.8	ug/l	154524	4	12.024	372103	50.0
Benzene, 1,2,4,...	12.920	10.6	ug/l	78995	4	12.024	372103	50.0
Benzene, 1,2,3,...	12.963	8.6	ug/l	64042	4	12.024	372103	50.0
1H-Indene, 2,3-...	13.292	12.9	ug/l	95593	4	12.024	372103	50.0
unknown13.420	13.420	5.6	ug/l	41502	4	12.024	372103	50.0