LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX113022\
Data File : VX033139.D
Acq On : 30 Nov 2022 18:01
Operator : JC/MD
Sample : N5815-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-03

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\82X112222W.M
Title : SW846 8260

Signal : TIC: VX033139.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.374	43	47	54	rBV	2015022	2345596	9.58%	1.087%
2	1.752	103	109	127	rVB	9904919	12849370	52.50%	5.956%
3	1.959	138	143	156	rVB	3104649	4509818	18.43%	2.090%
4	2.069	156	161	168	rBV	503386	722520	2.95%	0.335%
5	2.166	172	177	181	rVV	171265	252295	1.03%	0.117%
6	2.215	181	185	193	rVB	270410	422300	1.73%	0.196%
7	2.343	200	206	210	rBV	238920	437856	1.79%	0.203%
8	2.386	210	213	229	rVB	169064	337902	1.38%	0.157%
9	2.782	271	278	281	rVV	826511	1732047	7.08%	0.803%
10	2.831	281	286	290	rVV	1944802	4238268	17.32%	1.964%
11	2.867	290	292	308	rVB2	1015357	2020561	8.26%	0.937%
12	3.105	321	331	351	rVB	1439059	3201001	13.08%	1.484%
13	3.318	357	366	377	rBV2	169486	397309	1.62%	0.184%
14	3.446	377	387	400	rBV	669071	1481650	6.05%	0.687%
15	3.587	400	410	416	rBV2	140492	354907	1.45%	0.165%
16	3.672	416	424	428	rVV	137691	324110	1.32%	0.150%
17	3.733	428	434	443	rVB	239709	571384	2.33%	0.265%
18	3.837	443	451	458	rBV	128936	336178	1.37%	0.156%
19	3.952	466	470	483	rVB2	98421	284727	1.16%	0.132%
20	4.105	483	495	504	rVB2	167287	445661	1.82%	0.207%
21	4.221	504	514	519	rBV	180917	525211	2.15%	0.243%
22	4.312	519	529	543	rVB	2870361	8176279	33.40%	3.790%
23	4.562	561	570	590	rVB	114652	336282	1.37%	0.156%
24	5.196	666	674	690	rVB2	109047	364472	1.49%	0.169%
25	5.385	690	705	711	rBV2	90997	308014	1.26%	0.143%
26	5.483	711	721	737	rVV5	559291	2726896	11.14%	1.264%
27	5.623	737	744	764	rVB2	379743	1327344	5.42%	0.615%
28	5.830	765	778	791	rBV2	360978	1081756	4.42%	0.501%
29	6.043	804	813	822	rVV	426565	1149471	4.70%	0.533%
30	6.159	822	832	840	rVV3	258932	938800	3.84%	0.435%
31	6.257	840	848	857	rVV3	550838	1722463	7.04%	0.798%
32	6.361	857	865	876	rVB2	225693	646830	2.64%	0.300%
33	6.617	893	907	917	rBV5	240708	798576	3.26%	0.370%
34	6.763	924	931	936	rBV	221271	550245	2.25%	0.255%
35	6.848	939	945	959	rVB4	200284	630115	2.57%	0.292%
36	7.379	1017	1032	1040	rBV4	751027	2053633	8.39%	0.952%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX113022\
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Acq On : 30 Nov 2022 18:01
Operator : JC/MD
Sample : N5815-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-03

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\82X112222W.M
Title : SW846 8260

37	7.659	1071	1078	1090	rVB2	223048	499901	2.04%	0.232%
38	7.891	1104	1116	1123	rBV4	111836	353830	1.45%	0.164%
39	7.988	1123	1132	1143	rVB	243694	555178	2.27%	0.257%
40	8.110	1143	1152	1161	rBV2	367492	863933	3.53%	0.400%
41	8.439	1200	1206	1212	rBV3	157462	286214	1.17%	0.133%
42	8.506	1212	1217	1224	rVB3	175660	312804	1.28%	0.145%
43	8.610	1224	1234	1236	rBV4	123071	273890	1.12%	0.127%
44	8.647	1236	1240	1246	rVB	353058	587164	2.40%	0.272%
45	8.714	1246	1251	1257	rBV3	323780	558275	2.28%	0.259%
46	8.805	1261	1266	1270	rVV	180857	254404	1.04%	0.118%
47	10.055	1466	1471	1475	rVB	462734	626847	2.56%	0.291%
48	10.201	1488	1495	1501	rBV	17239664	24476357	100.00%	11.345%
49	10.311	1503	1513	1518	rVB	15610284	20781495	84.90%	9.632%
50	10.640	1562	1567	1579	rVB	558255	766344	3.13%	0.355%
51	10.963	1613	1620	1626	rBV	1991413	2551180	10.42%	1.182%
52	11.079	1634	1639	1644	rBV	326650	459852	1.88%	0.213%
53	11.305	1668	1676	1682	rBV	6176167	7491080	30.61%	3.472%
54	11.402	1682	1692	1696	rVV	8189310	9814671	40.10%	4.549%
55	11.451	1696	1700	1712	rVB	2104114	2608159	10.66%	1.209%
56	11.616	1722	1727	1734	rVB	3000374	3671909	15.00%	1.702%
57	11.756	1744	1750	1755	rBV	16394938	21029977	85.92%	9.748%
58	11.866	1761	1768	1770	rBV	299090	387865	1.58%	0.180%
59	11.890	1770	1772	1777	rVB	386415	426855	1.74%	0.198%
60	11.969	1781	1785	1789	rBV	242332	300419	1.23%	0.139%
61	12.024	1789	1794	1799	rVB2	592062	882352	3.60%	0.409%
62	12.091	1799	1805	1810	rBV	5088885	6203335	25.34%	2.875%
63	12.244	1824	1830	1837	rBV2	5994848	7847950	32.06%	3.638%
64	12.317	1837	1842	1850	rVB3	2298343	3742667	15.29%	1.735%
65	12.408	1852	1857	1861	rBV	322296	412358	1.68%	0.191%
66	12.463	1861	1866	1872	rVB	1274872	1508763	6.16%	0.699%
67	12.530	1872	1877	1880	rBV	1888760	2623793	10.72%	1.216%
68	12.555	1880	1881	1886	rVB	1311059	1117720	4.57%	0.518%
69	12.615	1886	1891	1894	rBV	3688841	4201665	17.17%	1.948%
70	12.646	1894	1896	1900	rVB	520394	531641	2.17%	0.246%
71	12.701	1900	1905	1913	rBV2	1456872	2207247	9.02%	1.023%
72	12.798	1916	1921	1925	rBV	297890	374334	1.53%	0.174%
73	12.847	1925	1929	1934	rVB	901562	1086573	4.44%	0.504%
74	12.920	1934	1941	1944	rBV	1663384	2179237	8.90%	1.010%
75	12.963	1944	1948	1954	rVB	1878460	2445812	9.99%	1.134%
76	13.024	1954	1958	1960	rBV	204910	258626	1.06%	0.120%

LSC Area Percent Report

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Sample : N5815-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-03

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

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Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Title : SW846 8260

77	13.170	1978	1982	1986	rVB	1312478	1529445	6.25%	0.709%
78	13.256	1992	1996	1998	rBV2	220139	307485	1.26%	0.143%
79	13.292	1998	2002	2008	rVV2	2749000	3722403	15.21%	1.725%
80	13.420	2020	2023	2032	rVB	280746	397623	1.62%	0.184%
81	13.579	2046	2049	2057	rVV5	365455	826823	3.38%	0.383%
82	13.664	2060	2063	2068	rVB2	199775	346923	1.42%	0.161%
83	13.780	2076	2082	2086	rBV	3674324	4895321	20.00%	2.269%
84	13.823	2086	2089	2094	rVB2	804587	967570	3.95%	0.448%
85	14.213	2149	2153	2159	rBV	156618	249531	1.02%	0.116%
86	14.640	2218	2223	2228	rVB	2268041	2789760	11.40%	1.293%
87	14.780	2241	2246	2255	rVB	1256741	1548590	6.33%	0.718%

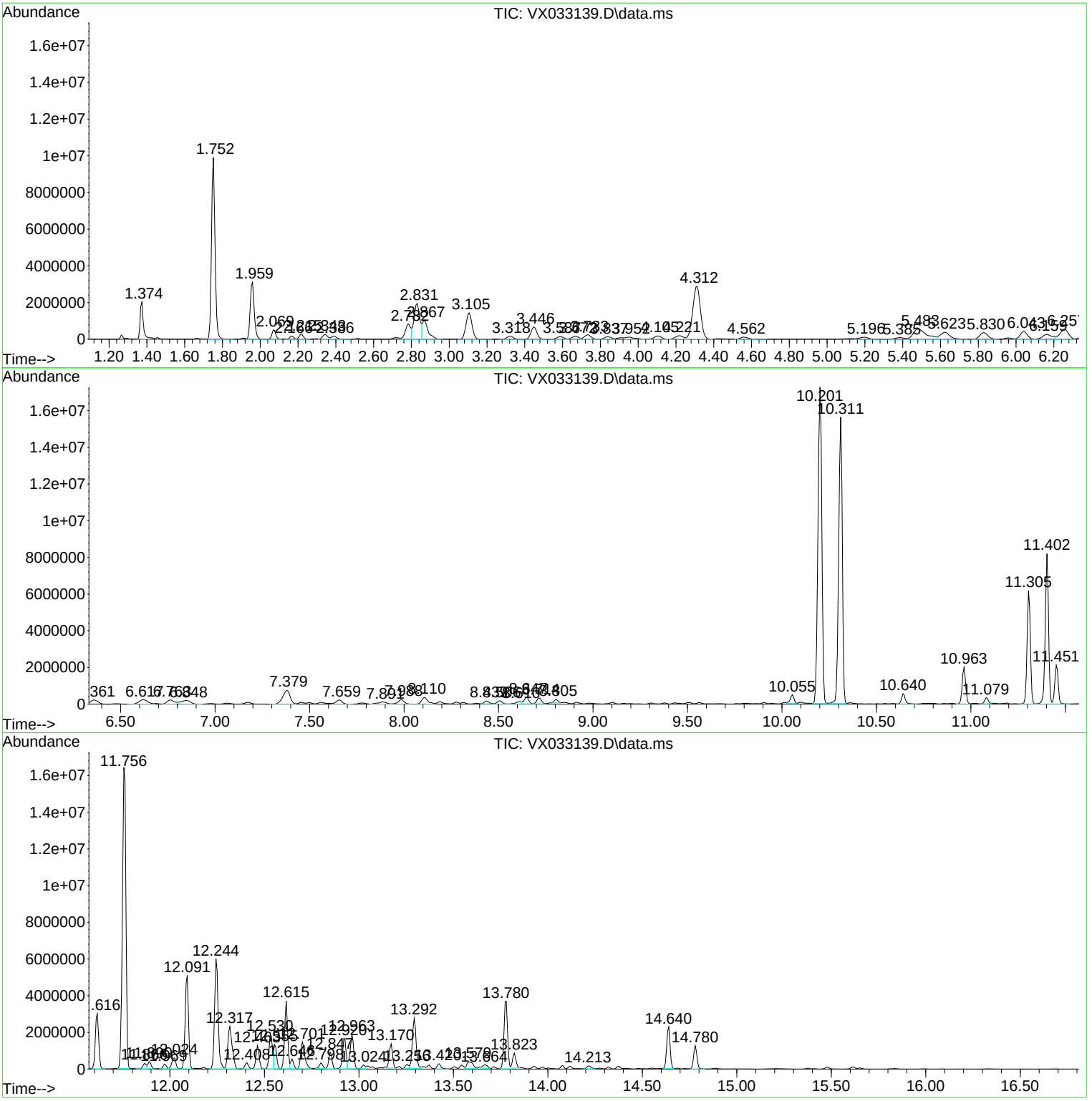
Sum of corrected areas: 215745997

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX113022\
Data File : VX033139.D
Acq On : 30 Nov 2022 18:01
Operator : JC/MD
Sample : N5815-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-03

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX113022\
Data File : VX033139.D
Acq On : 30 Nov 2022 18:01
Operator : JC/MD
Sample : N5815-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-03

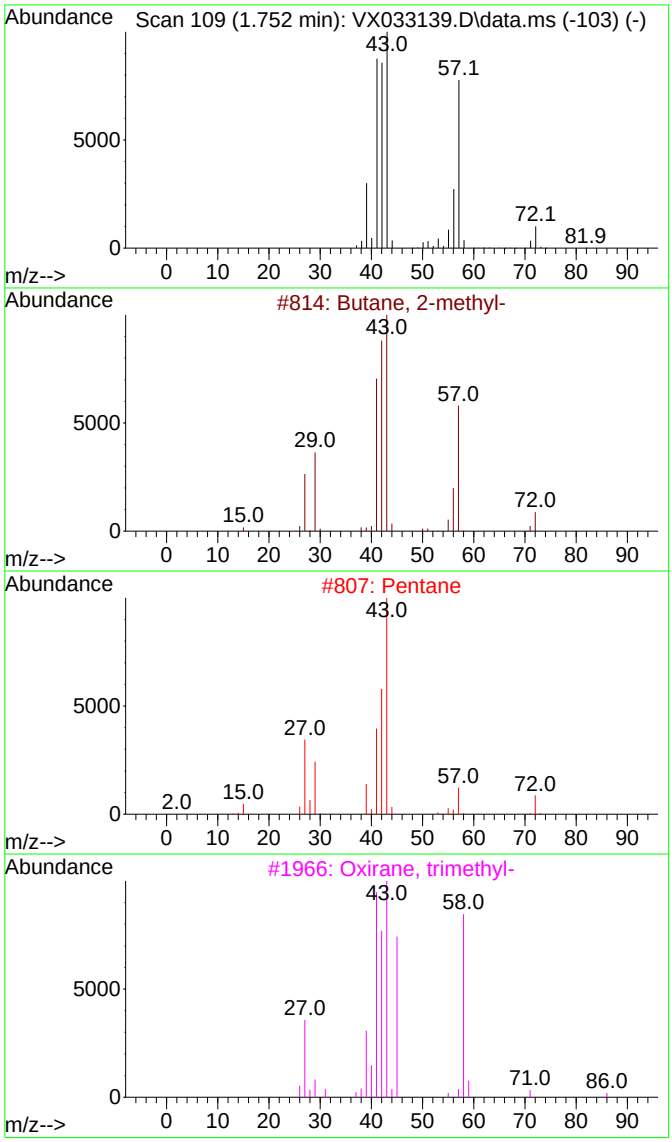
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

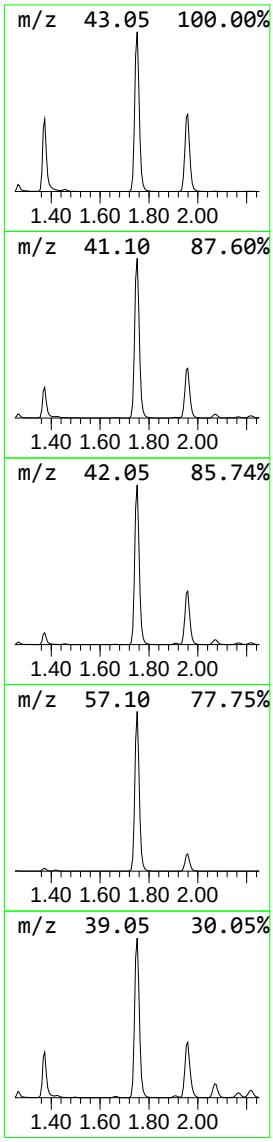
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.752	484.03 ug/l	12849400	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	91
2	Pentane	72	C5H12	000109-66-0	42
3	Oxirane, trimethyl-	86	C5H10O	005076-19-7	33
4	3-Buten-1-ol	72	C4H8O	000627-27-0	25
5	Propane, 2-methyl-2-nitro-	103	C4H9NO2	000594-70-7	10





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX113022\
Data File : VX033139.D
Acq On : 30 Nov 2022 18:01
Operator : JC/MD
Sample : N5815-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-03

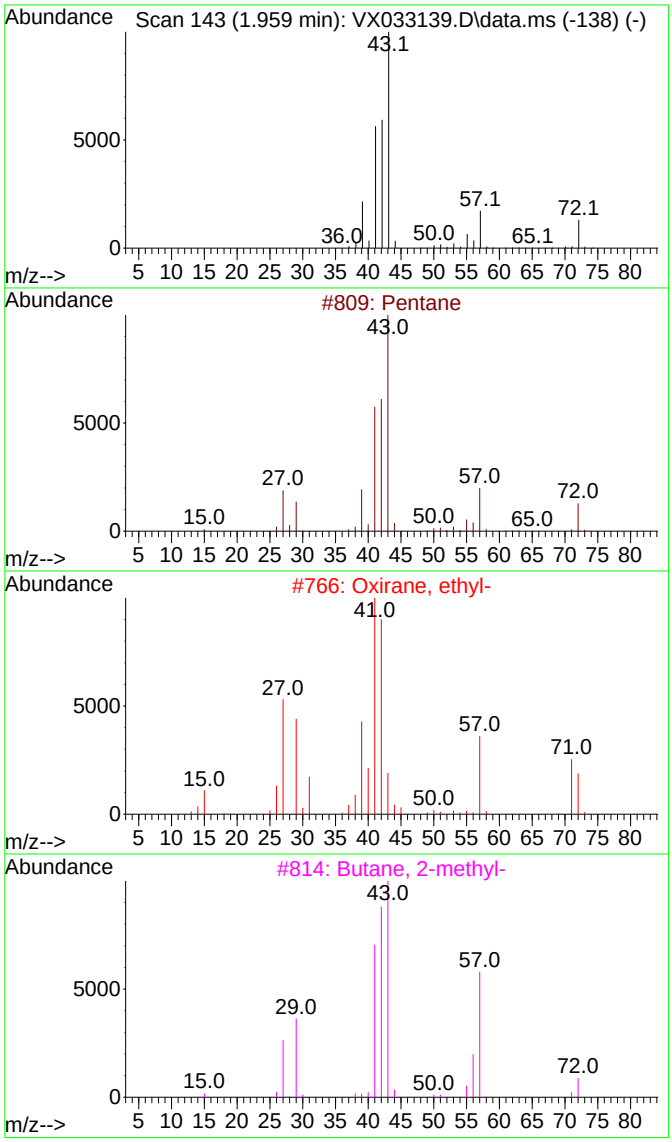
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

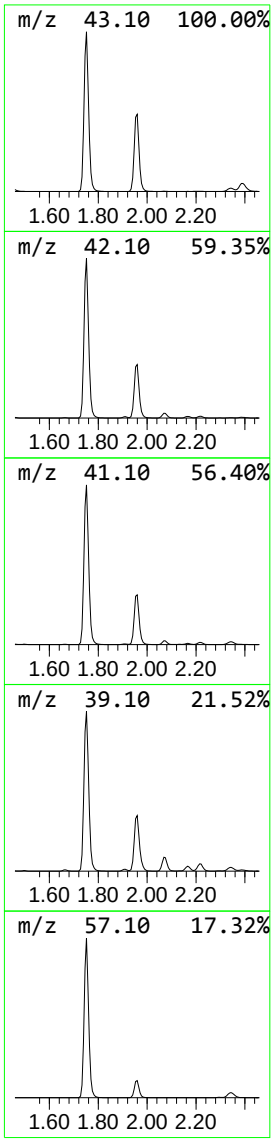
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Pentane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.959	169.88 ug/l	4509820	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane	72	C5H12	000109-66-0	91
2	Oxirane, ethyl-	72	C4H8O	000106-88-7	50
3	Butane, 2-methyl-	72	C5H12	000078-78-4	38
4	2(3H)-Furanone, dihydro-4-methyl-	100	C5H8O2	001679-49-8	25
5	2-Propanone, hydrazone	72	C3H8N2	005281-20-9	10





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX113022\
Data File : VX033139.D
Acq On : 30 Nov 2022 18:01
Operator : JC/MD
Sample : N5815-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-03

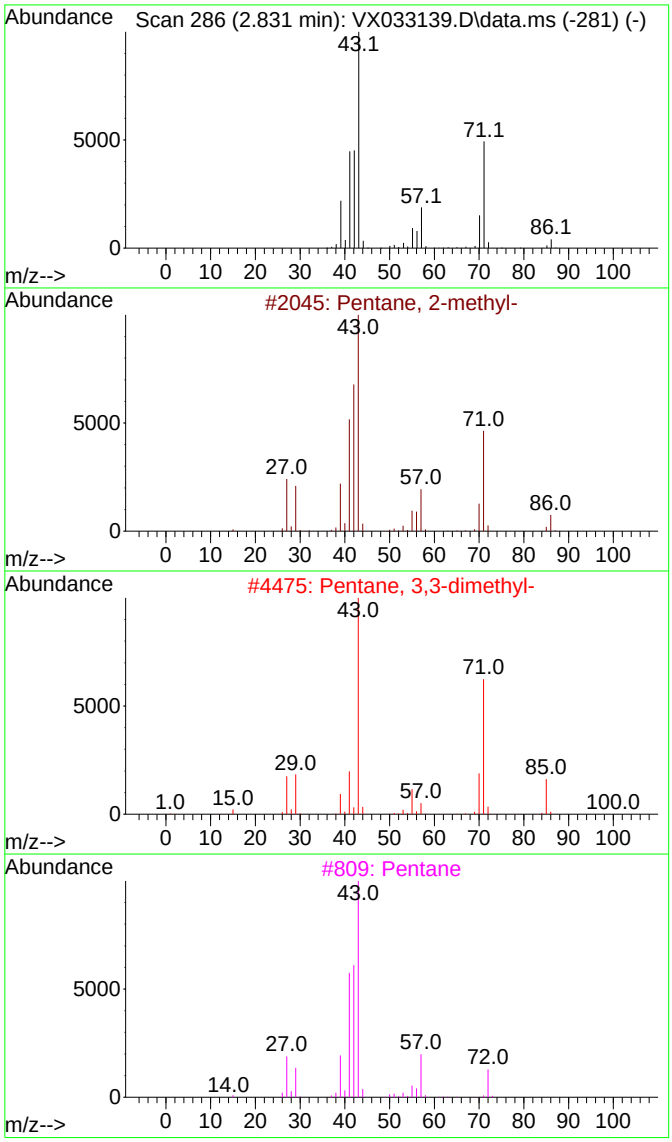
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

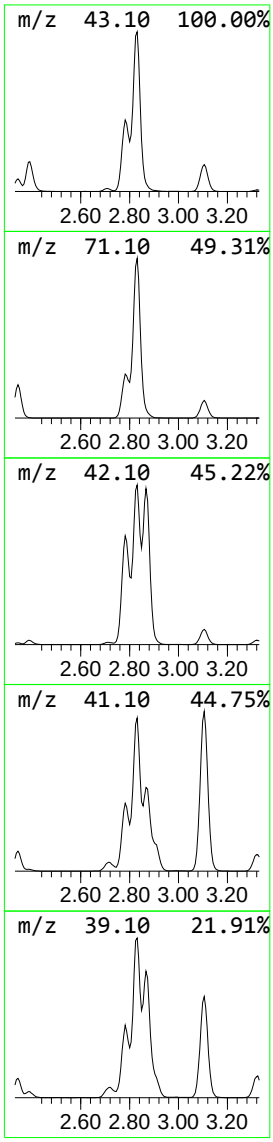
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Pentane, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.831	159.65 ug/l	4238270	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2-methyl-	86	C6H14	000107-83-5	94
2	Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	47
3	Pentane	72	C5H12	000109-66-0	46
4	1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	45
5	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	43





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX113022\
Data File : VX033139.D
Acq On : 30 Nov 2022 18:01
Operator : JC/MD
Sample : N5815-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-03

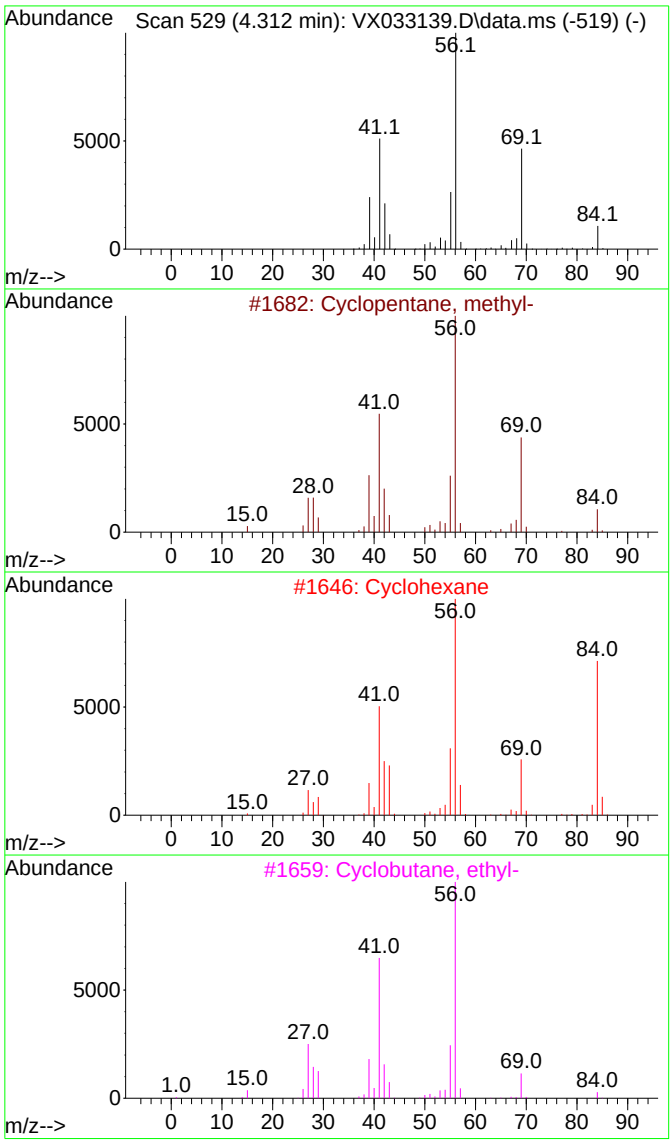
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

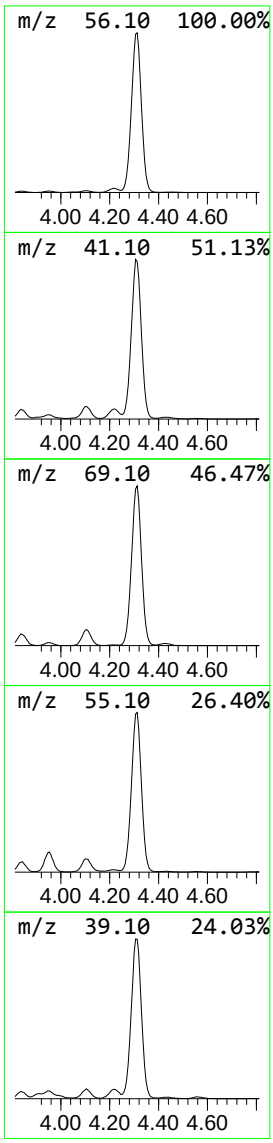
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.312	307.99 ug/l	8176280	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2	Cyclohexane	84	C6H12	000110-82-7	83
3	Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
4	Cyclopropane, propyl-	84	C6H12	002415-72-7	59
5	Cyclobutane	56	C4H8	000287-23-0	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX113022\
Data File : VX033139.D
Acq On : 30 Nov 2022 18:01
Operator : JC/MD
Sample : N5815-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-03

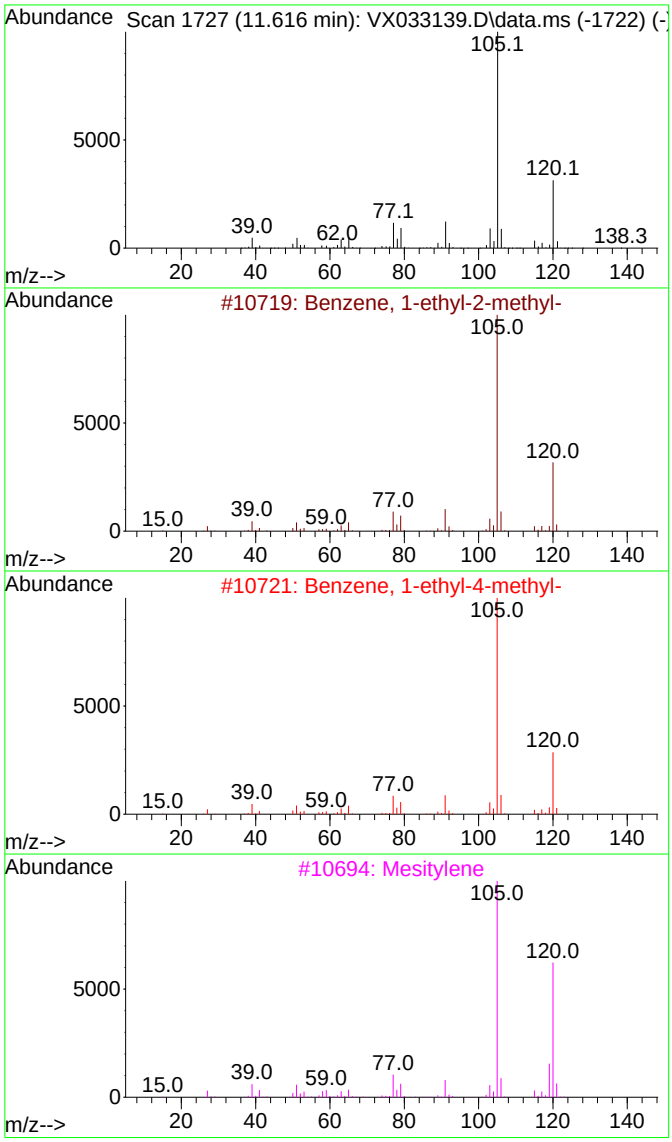
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

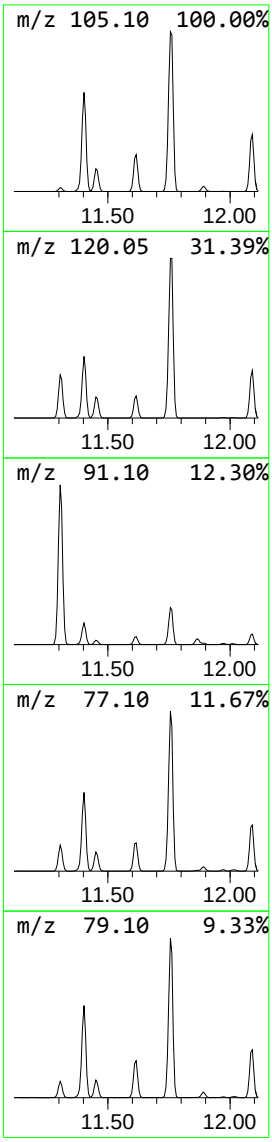
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1-ethyl-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.616	208.07 ug/l	3671910	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX113022\
Data File : VX033139.D
Acq On : 30 Nov 2022 18:01
Operator : JC/MD
Sample : N5815-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-03

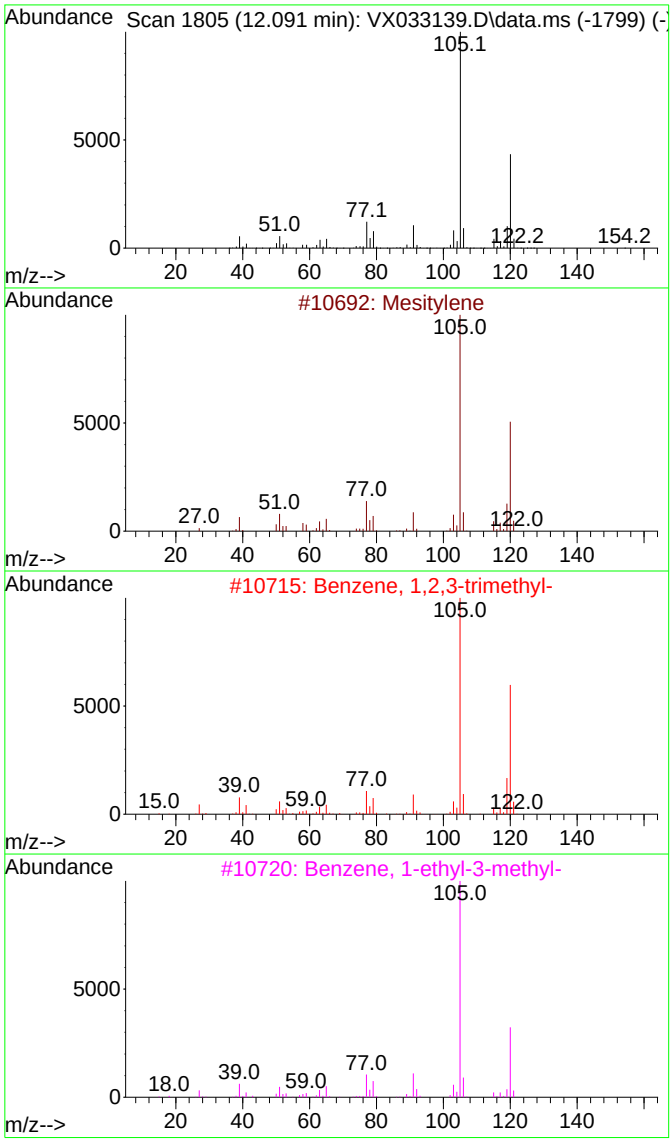
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

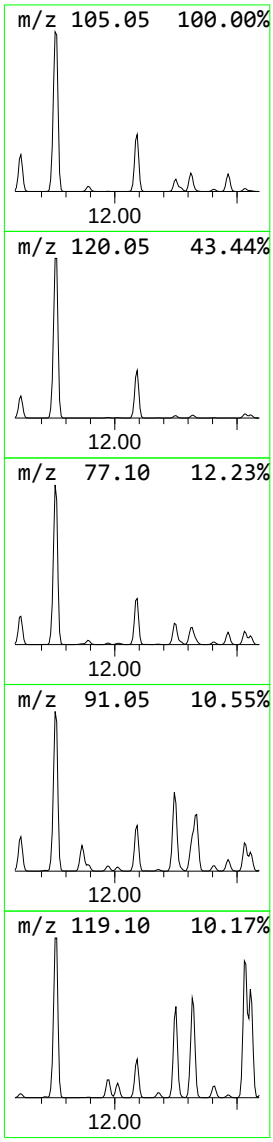
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.091	351.52 ug/l	6203340	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	97
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX113022\
Data File : VX033139.D
Acq On : 30 Nov 2022 18:01
Operator : JC/MD
Sample : N5815-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-03

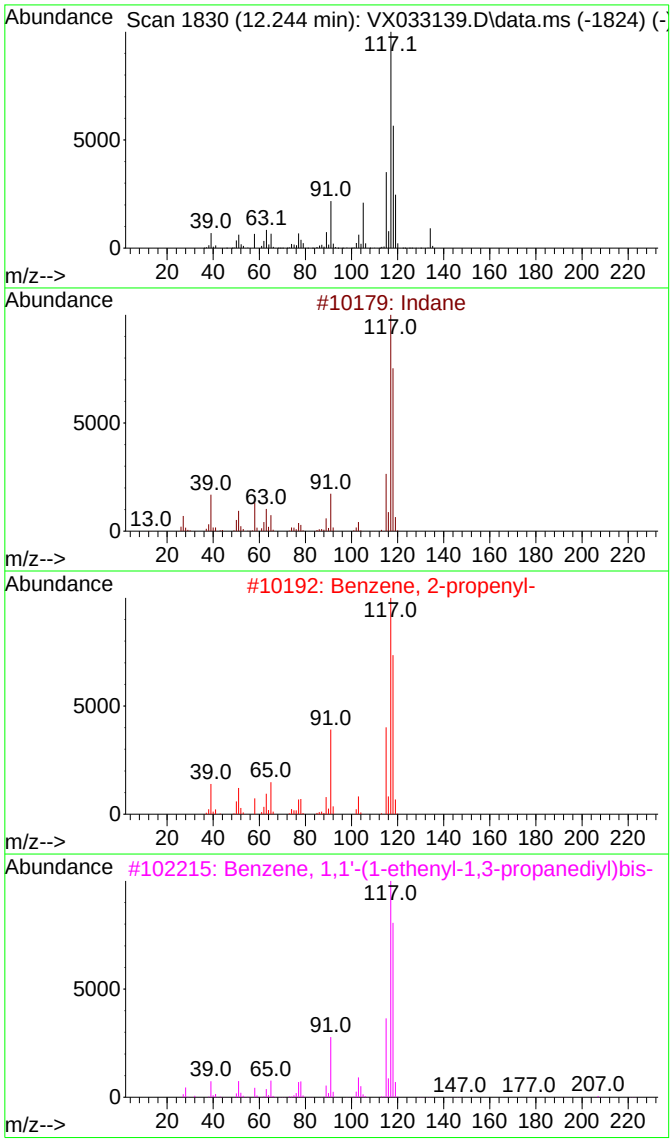
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

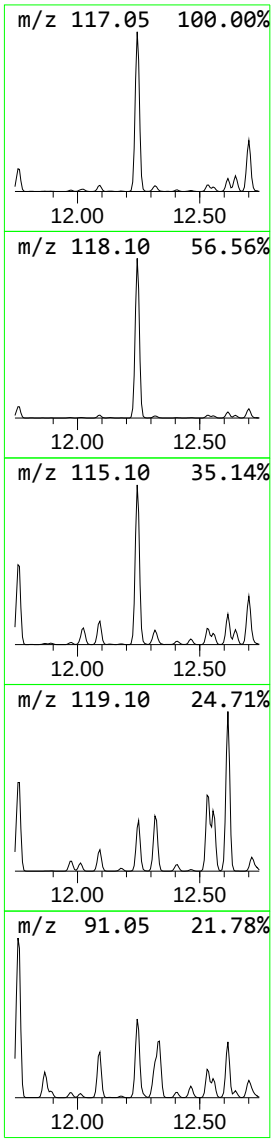
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	444.72 ug/l	7847950	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	64
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
3	Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	64
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	46





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX113022\
Data File : VX033139.D
Acq On : 30 Nov 2022 18:01
Operator : JC/MD
Sample : N5815-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-03

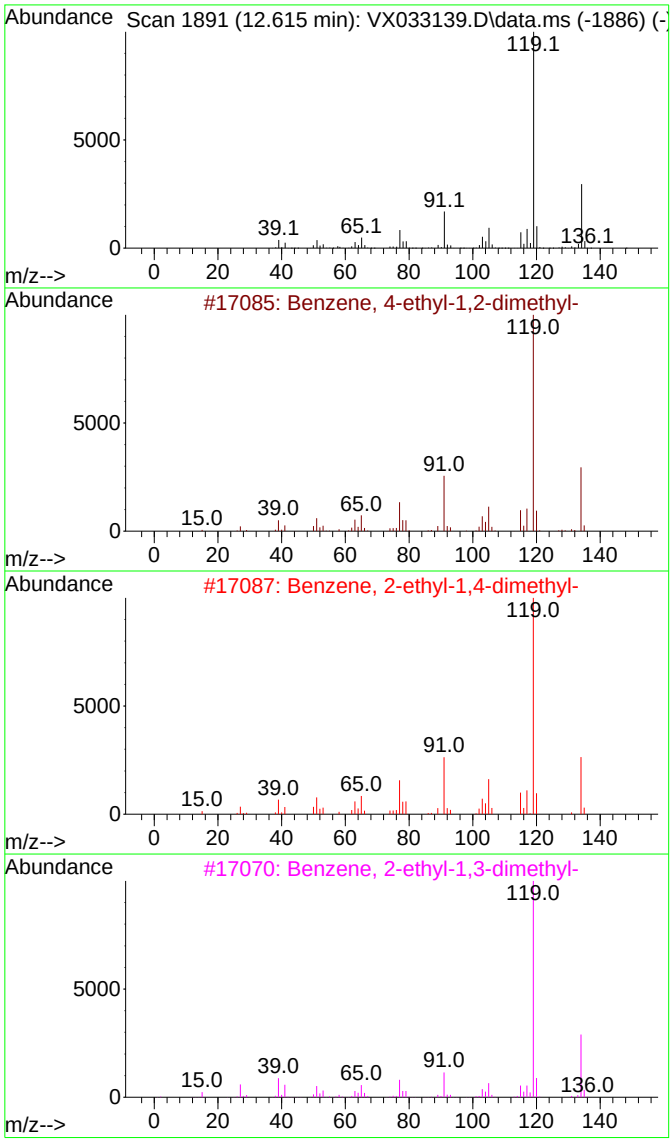
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

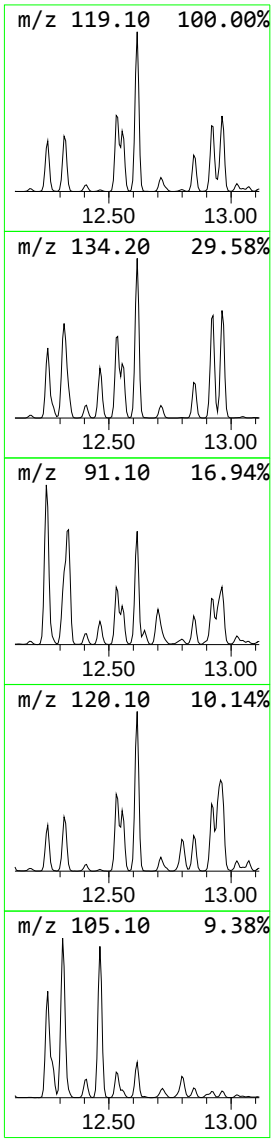
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.616	238.10 ug/l	4201670	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5	o-Cymene	134	C10H14	000527-84-4	95





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX113022\
Data File : VX033139.D
Acq On : 30 Nov 2022 18:01
Operator : JC/MD
Sample : N5815-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-03

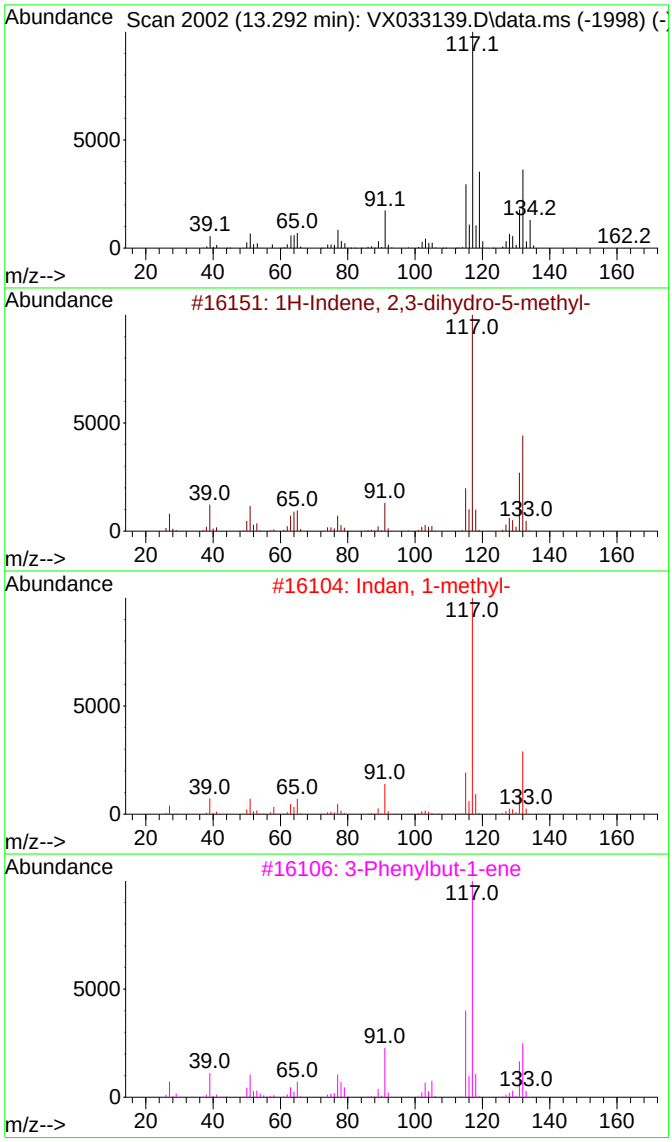
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

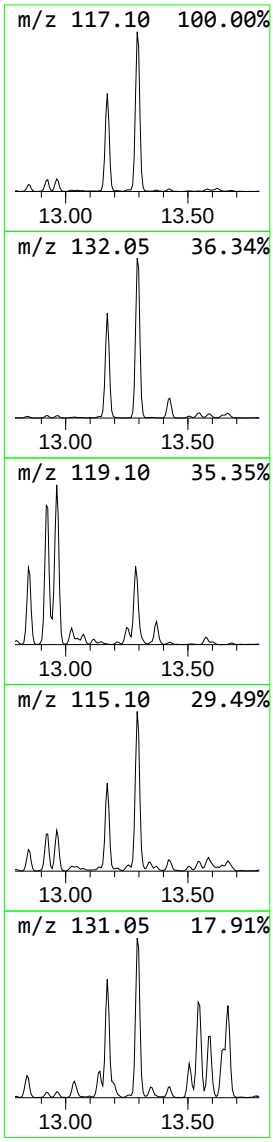
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	210.94 ug/l	3722400	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	89
2	Indan, 1-methyl-	132	C10H12	000767-58-8	81
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
4	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	81
5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	76





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX113022\
Data File : VX033139.D
Acq On : 30 Nov 2022 18:01
Operator : JC/MD
Sample : N5815-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-03

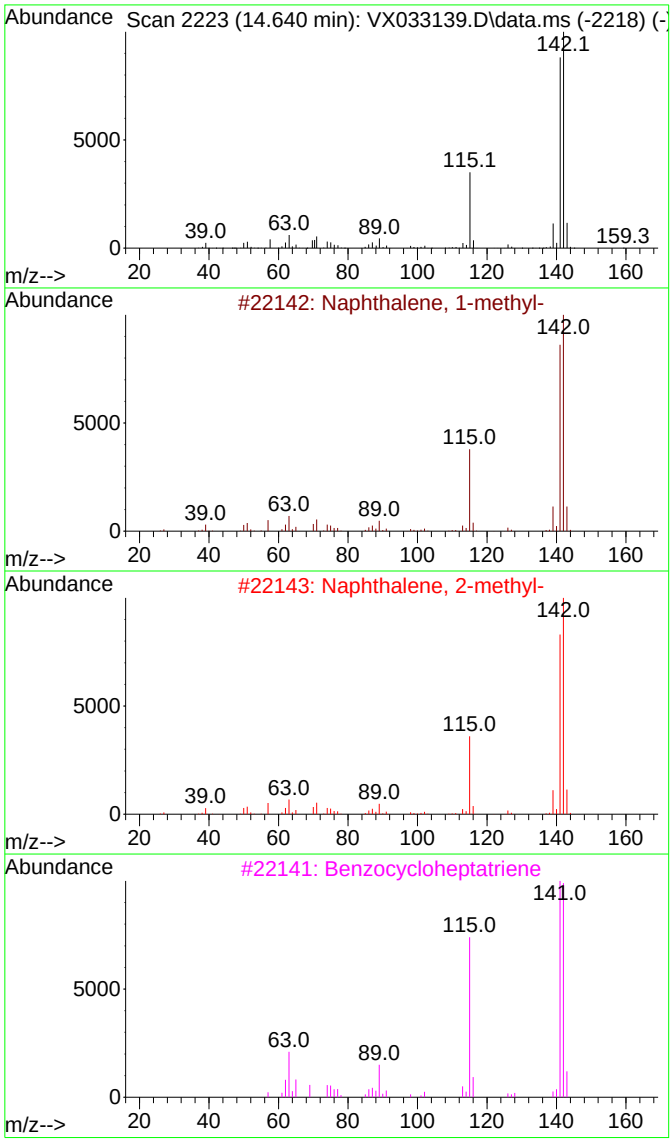
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

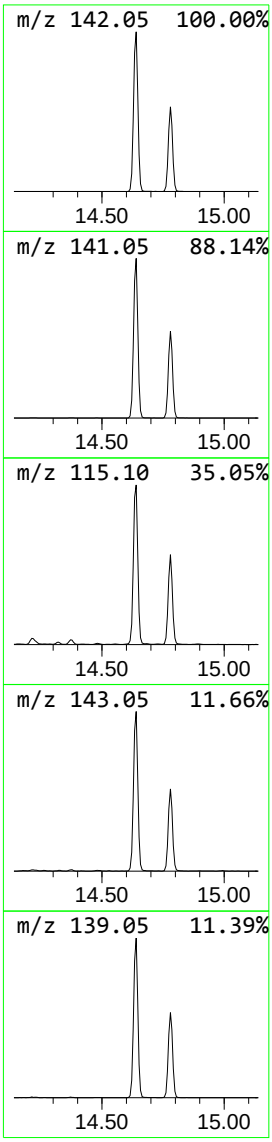
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.640	158.09 ug/l	2789760	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3	Benzocycloheptatriene	142	C11H10	000264-09-5	91
4	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX113022\
Data File : VX033139.D
Acq On : 30 Nov 2022 18:01
Operator : JC/MD
Sample : N5815-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-03

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.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	484.0	ug/l	12849400	1	5.556	1327340	50.0
Pentane	1.959	169.9	ug/l	4509820	1	5.556	1327340	50.0
Pentane, 2-methyl-	2.831	159.7	ug/l	4238270	1	5.556	1327340	50.0
Cyclopentane, m...	4.312	308.0	ug/l	8176280	1	5.556	1327340	50.0
Benzene, 1-ethy...	11.616	208.1	ug/l	3671910	4	12.024	882352	50.0
Benzene, 1,2,3-...	12.091	351.5	ug/l	6203340	4	12.024	882352	50.0
Indane	12.244	444.7	ug/l	7847950	4	12.024	882352	50.0
Benzene, 4-ethy...	12.616	238.1	ug/l	4201670	4	12.024	882352	50.0
1H-Indene, 2,3-...	13.292	210.9	ug/l	3722400	4	12.024	882352	50.0
Naphthalene, 1-...	14.640	158.1	ug/l	2789760	4	12.024	882352	50.0