LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100322\
Data File : VX031845.D
Acq On : 03 Oct 2022 19:48
Operator : JC/MD
Sample : N4860-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2R

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\82X100322W.M
Title : SW846 8260

Signal : TIC: VX031845.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.282	26	32	51	rBV	55805	234239	37.14%	2.407%
2	5.385	694	705	722	rBV	79006	232148	36.81%	2.385%
3	5.549	722	732	749	rVB2	77880	220287	34.93%	2.264%
4	5.958	789	799	810	rBV	77120	208176	33.01%	2.139%
5	6.763	919	931	945	rBV	110523	260843	41.36%	2.680%
6	8.646	1233	1240	1252	rBV	379213	609592	96.65%	6.264%
7	10.055	1464	1471	1480	rBV	272763	359973	57.07%	3.699%
8	10.274	1503	1507	1517	rVB5	4255	8894	1.41%	0.091%
9	10.585	1550	1558	1565	rBV8	3556	9227	1.46%	0.095%
10	10.963	1616	1620	1624	rBV	6885	8976	1.42%	0.092%
11	11.079	1634	1639	1645	rBV	301530	393645	62.41%	4.045%
12	11.219	1659	1662	1667	rVB4	6068	8533	1.35%	0.088%
13	11.615	1721	1727	1735	rBV3	7554	14085	2.23%	0.145%
14	11.713	1739	1743	1746	rBV	6600	8383	1.33%	0.086%
15	11.853	1759	1766	1769	rBV4	4151	9211	1.46%	0.095%
16	11.890	1769	1772	1777	rVB	11501	15242	2.42%	0.157%
17	12.024	1788	1794	1800	rBV	345797	419344	66.49%	4.309%
18	12.176	1813	1819	1825	rVB	25403	35143	5.57%	0.361%
19	12.243	1825	1830	1836	rBV2	515004	630725	100.00%	6.481%
20	12.310	1837	1841	1849	rVB2	33582	50375	7.99%	0.518%
21	12.402	1849	1856	1864	rBV2	22371	35213	5.58%	0.362%
22	12.463	1864	1866	1871	rVB5	3704	6792	1.08%	0.070%
23	12.585	1880	1886	1887	rBV6	5418	7517	1.19%	0.077%
24	12.609	1887	1890	1893	rVV	18481	24655	3.91%	0.253%
25	12.646	1893	1896	1900	rVV	93965	108269	17.17%	1.113%
26	12.701	1900	1905	1912	rVB	218617	301965	47.88%	3.103%
27	12.762	1912	1915	1920	rVB2	11561	13006	2.06%	0.134%
28	12.841	1920	1928	1931	rVB2	31941	58298	9.24%	0.599%
29	12.920	1933	1941	1946	rBV2	29570	58905	9.34%	0.605%
30	12.981	1946	1951	1955	rVB3	11867	16466	2.61%	0.169%
31	13.030	1955	1959	1965	rVB2	53101	78093	12.38%	0.802%
32	13.097	1965	1970	1973	rBV	21203	32608	5.17%	0.335%
33	13.140	1973	1977	1980	rBV	48645	57913	9.18%	0.595%
34	13.170	1980	1982	1983	rBV2	10763	9498	1.51%	0.098%
35	13.194	1983	1986	1991	rVB	37686	47883	7.59%	0.492%
36	13.261	1991	1997	1998	rBV2	15418	24655	3.91%	0.253%

LSC Area Percent Report

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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2R

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\82X100322W.M
Title : SW846 8260

37	13.292	1998	2002	2008	rVB	328376	403162	63.92%	4.143%
38	13.347	2008	2011	2014	rBV2	28020	36557	5.80%	0.376%
39	13.389	2015	2018	2020	rBV3	11685	13974	2.22%	0.144%
40	13.420	2020	2023	2032	rVB3	42903	71073	11.27%	0.730%
41	13.505	2033	2037	2040	rBV2	21389	33262	5.27%	0.342%
42	13.542	2040	2043	2046	rBV	62925	78338	12.42%	0.805%
43	13.585	2046	2050	2053	rBV2	119752	140353	22.25%	1.442%
44	13.664	2060	2063	2068	rVB	215430	277605	44.01%	2.852%
45	13.749	2072	2077	2081	rBV2	25323	37605	5.96%	0.386%
46	13.798	2081	2085	2090	rBV3	28745	42650	6.76%	0.438%
47	13.853	2090	2094	2099	rVB2	131650	173479	27.50%	1.783%
48	13.926	2099	2106	2110	rBV	228666	298432	47.32%	3.067%
49	13.969	2110	2113	2117	rVV	86637	114966	18.23%	1.181%
50	14.005	2117	2119	2125	rVB	27756	33300	5.28%	0.342%
51	14.072	2125	2130	2134	rBV2	84386	117405	18.61%	1.206%
52	14.103	2134	2135	2138	rVV2	14034	12135	1.92%	0.125%
53	14.139	2138	2141	2143	rVV3	11026	16338	2.59%	0.168%
54	14.164	2143	2145	2148	rVV2	19316	20733	3.29%	0.213%
55	14.213	2148	2153	2163	rVV	185071	297420	47.16%	3.056%
56	14.286	2163	2165	2167	rVV	25883	24874	3.94%	0.256%
57	14.322	2167	2171	2175	rVV	126304	157496	24.97%	1.618%
58	14.371	2175	2179	2184	rVV	162151	215414	34.15%	2.213%
59	14.420	2184	2187	2192	rVB	45292	60275	9.56%	0.619%
60	14.481	2192	2197	2202	rBV2	265503	358287	56.81%	3.682%
61	14.523	2202	2204	2210	rVV3	23774	35108	5.57%	0.361%
62	14.615	2210	2219	2225	rVV4	152132	317735	50.38%	3.265%
63	14.676	2225	2229	2233	rVB	95106	136211	21.60%	1.400%
64	14.725	2233	2237	2243	rBV3	46251	78357	12.42%	0.805%
65	14.786	2243	2247	2250	rVV2	65381	123763	19.62%	1.272%
66	14.816	2250	2252	2255	rVV2	78801	104410	16.55%	1.073%
67	14.847	2255	2257	2261	rVV2	57552	70783	11.22%	0.727%
68	14.907	2261	2267	2273	rVV5	49666	125369	19.88%	1.288%
69	14.962	2273	2276	2283	rVB5	29128	68754	10.90%	0.706%
70	15.048	2285	2290	2294	rBV2	31331	48473	7.69%	0.498%
71	15.182	2306	2312	2316	rBV2	138101	243747	38.65%	2.505%
72	15.249	2320	2323	2331	rVB6	27038	61377	9.73%	0.631%
73	15.322	2331	2335	2339	rBV5	12153	18281	2.90%	0.188%
74	15.371	2339	2343	2346	rBV3	25732	39168	6.21%	0.402%
75	15.414	2346	2350	2353	rVB2	21832	32650	5.18%	0.335%
76	15.481	2358	2361	2364	rVV2	22240	28752	4.56%	0.295%

LSC Area Percent Report

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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2R

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	15.523	2365	2368	2373	rVB4	26683	48459	7.68%	0.498%
78	15.615	2379	2383	2387	rVB	31728	42785	6.78%	0.440%
79	15.743	2398	2404	2408	rBV7	17372	40487	6.42%	0.416%
80	15.810	2411	2415	2418	rVV	50282	80891	12.83%	0.831%
81	15.846	2418	2421	2430	rVB3	44132	75353	11.95%	0.774%
82	15.987	2440	2444	2448	rBV2	25334	38594	6.12%	0.397%
83	16.084	2457	2460	2465	rBV7	9044	15554	2.47%	0.160%
84	16.261	2485	2489	2492	rBV2	14467	26406	4.19%	0.271%
85	16.377	2506	2508	2512	rBV3	4696	6783	1.08%	0.070%
86	16.426	2512	2516	2520	rVB	12746	18087	2.87%	0.186%
87	16.474	2521	2524	2528	rBV5	8113	12094	1.92%	0.124%
88	16.529	2529	2533	2539	rVV6	14961	36432	5.78%	0.374%
89	16.596	2540	2544	2549	rVV2	27516	50152	7.95%	0.515%
90	16.657	2549	2554	2561	rVB3	30058	53080	8.42%	0.545%

Sum of corrected areas: 9732005

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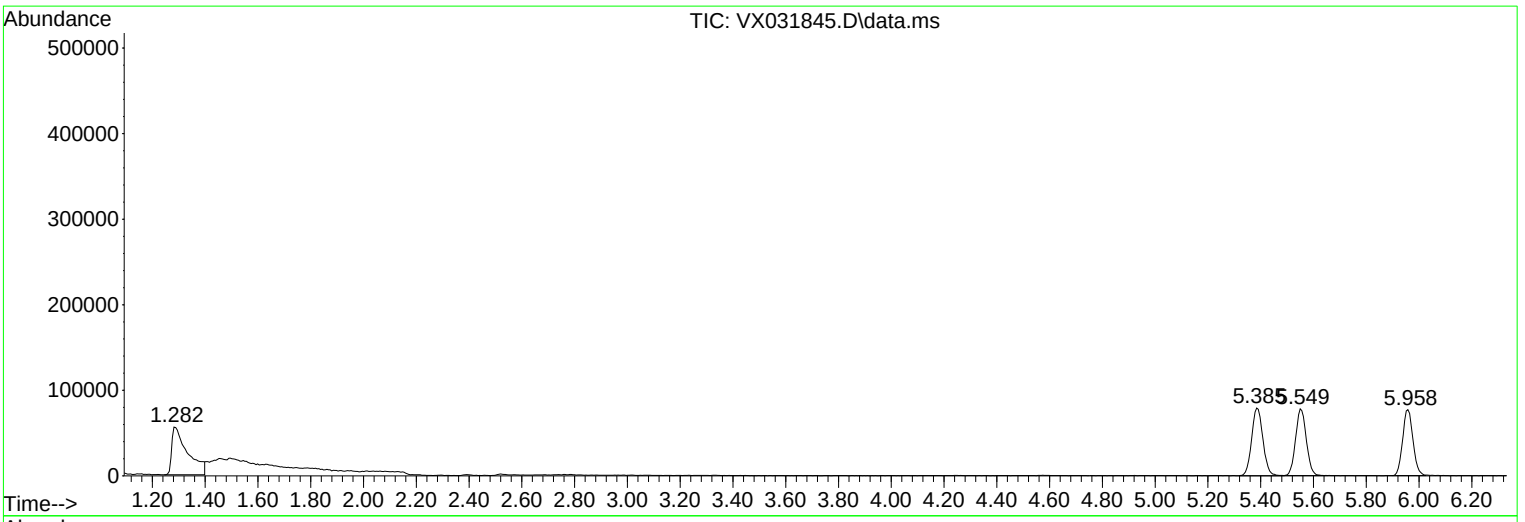
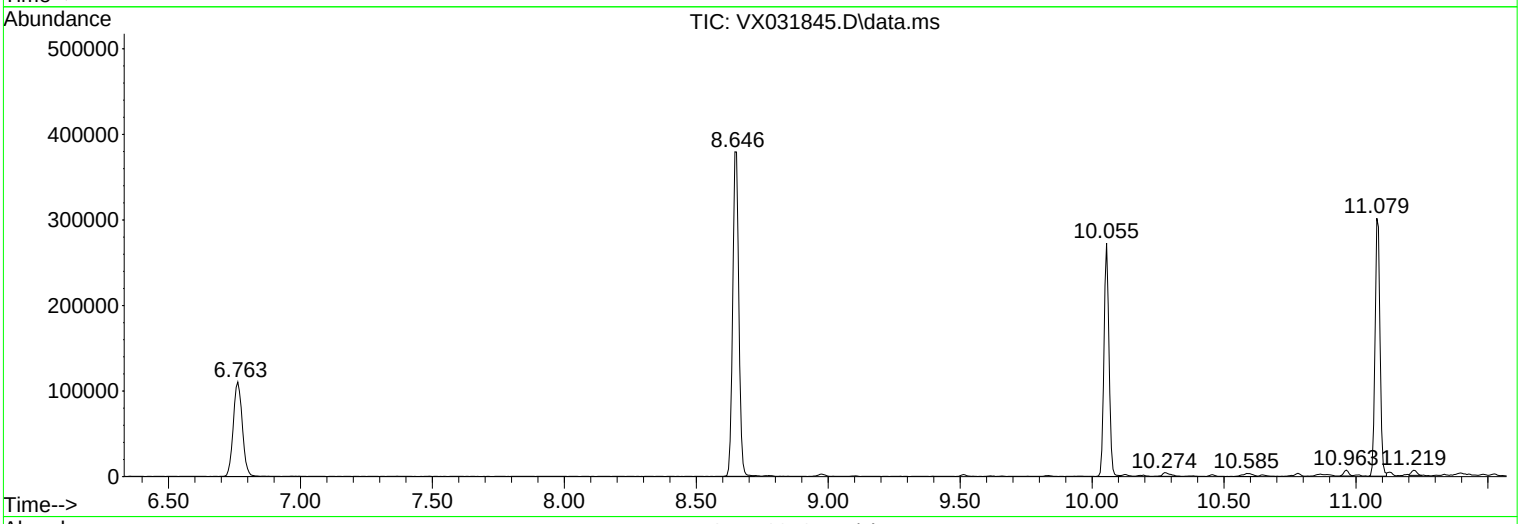
ALS Vial : 23 Sample Multiplier: 1

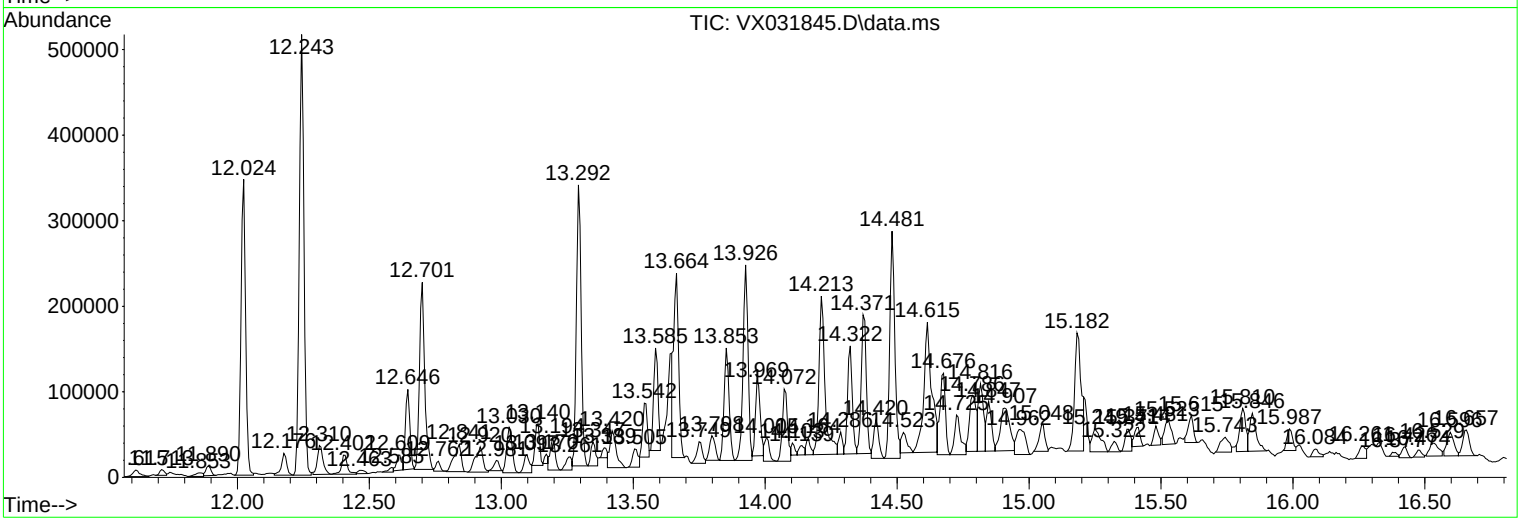
Instrument :
MSVOA_X
ClientSampleId :
MW-2R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100322W.M
Quant Title : SW846 8260

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







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Data File : VX031845.D
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Operator : JC/MD
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Instrument :
MSVOA_X
ClientSampleId :
MW-2R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100322W.M
Quant Title : SW846 8260

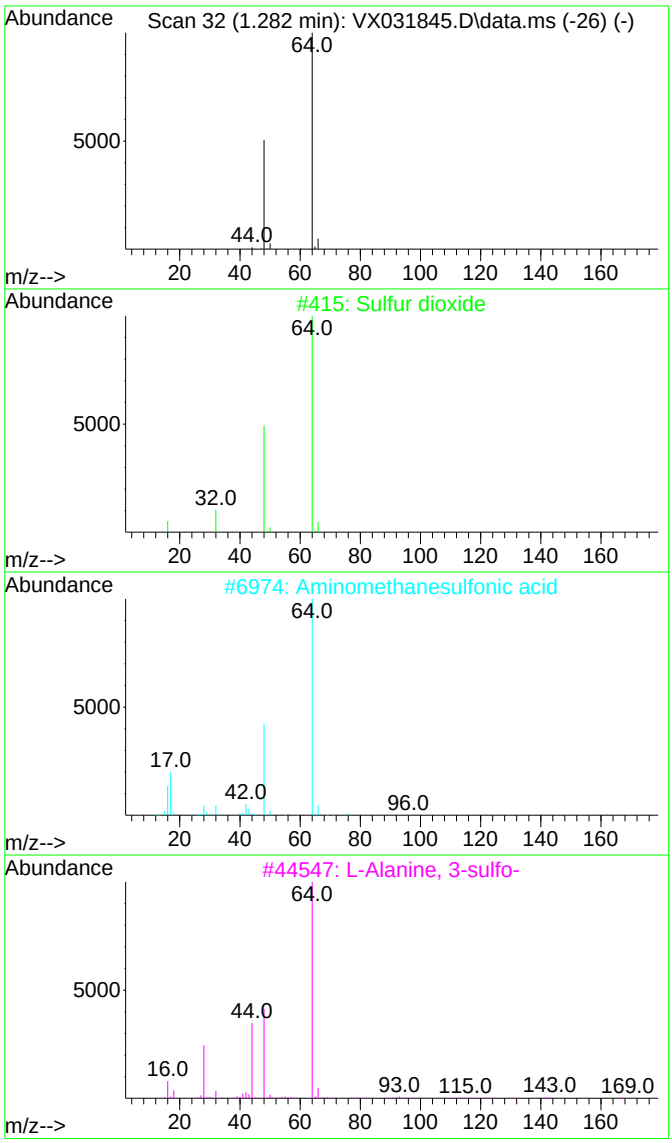
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

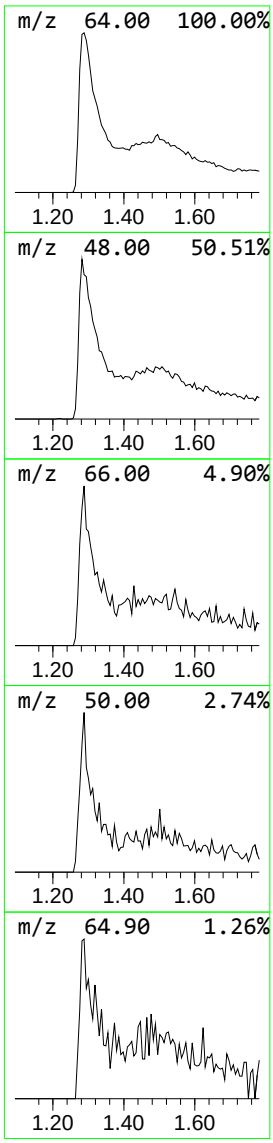
****

Peak Number 1 Sulfur dioxide Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.282	53.17 ug/l	234239	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\VX100322\
Data File : VX031845.D
Acq On : 03 Oct 2022 19:48
Operator : JC/MD
Sample : N4860-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2R

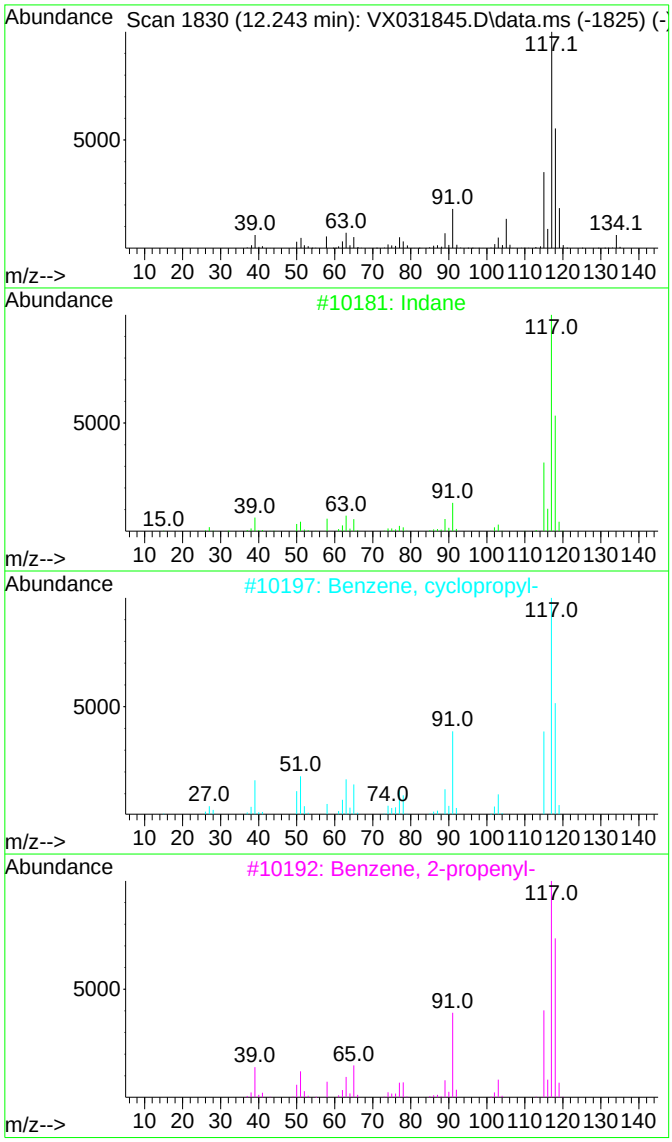
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100322W.M
Quant Title : SW846 8260

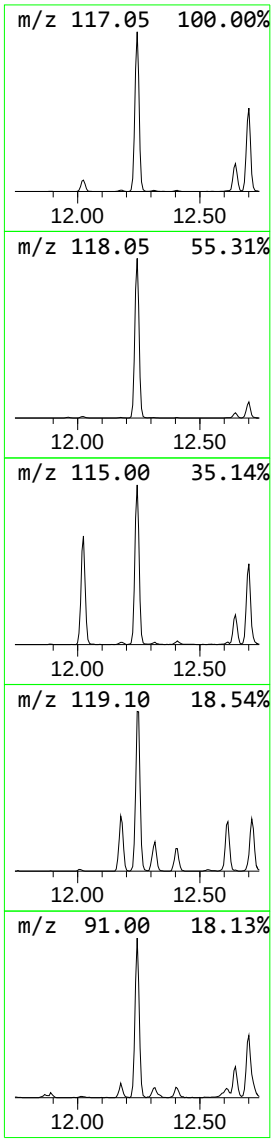
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.243	75.20 ug/l	630725	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	93
2	Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	68
4	Deltacyclene	118	C9H10	007785-10-6	58
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	53





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Instrument :
MSVOA_X
ClientSampleId :
MW-2R

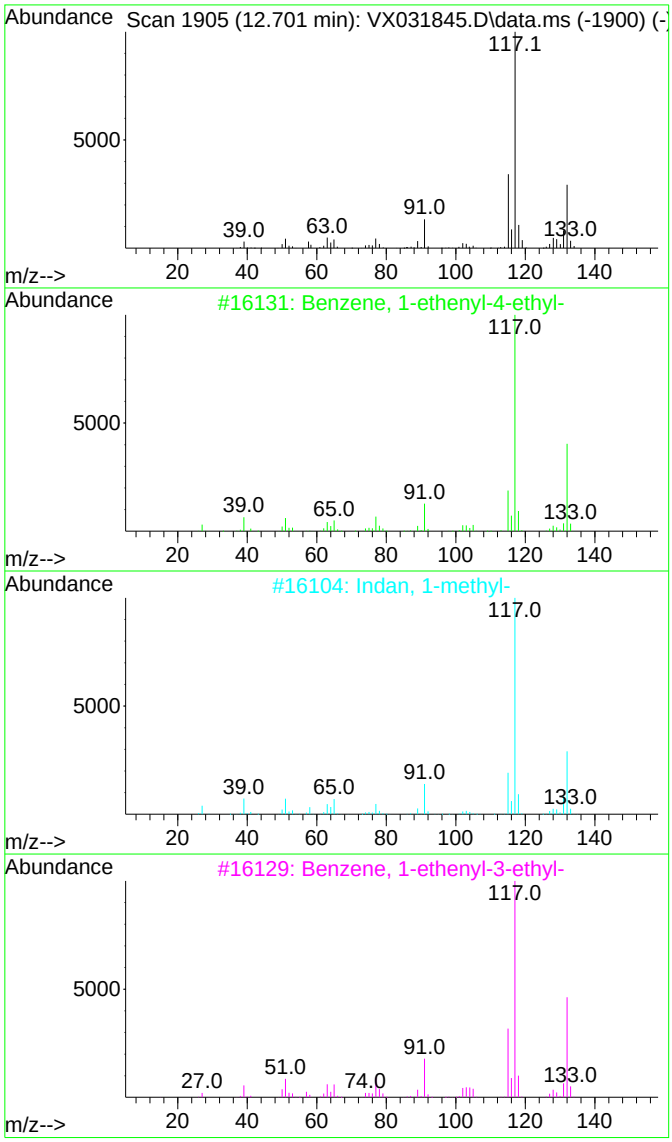
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100322W.M
Quant Title : SW846 8260

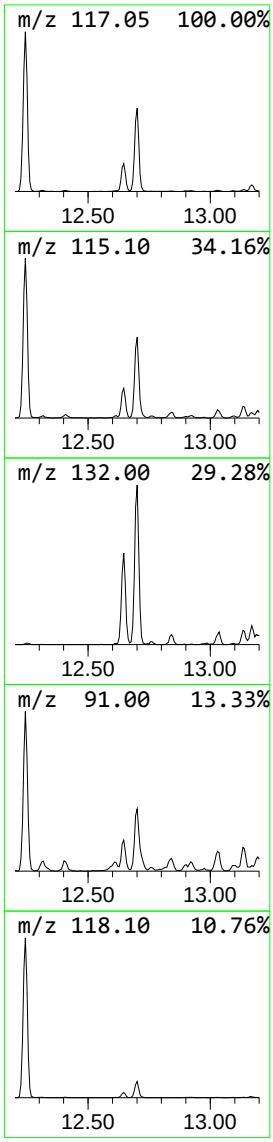
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	36.00 ug/l	301965	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
2	Indan, 1-methyl-	132	C10H12	000767-58-8	87
3	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	83
5	1-Phenyl-1-butene	132	C10H12	000824-90-8	83





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100322\
Data File : VX031845.D
Acq On : 03 Oct 2022 19:48
Operator : JC/MD
Sample : N4860-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2R

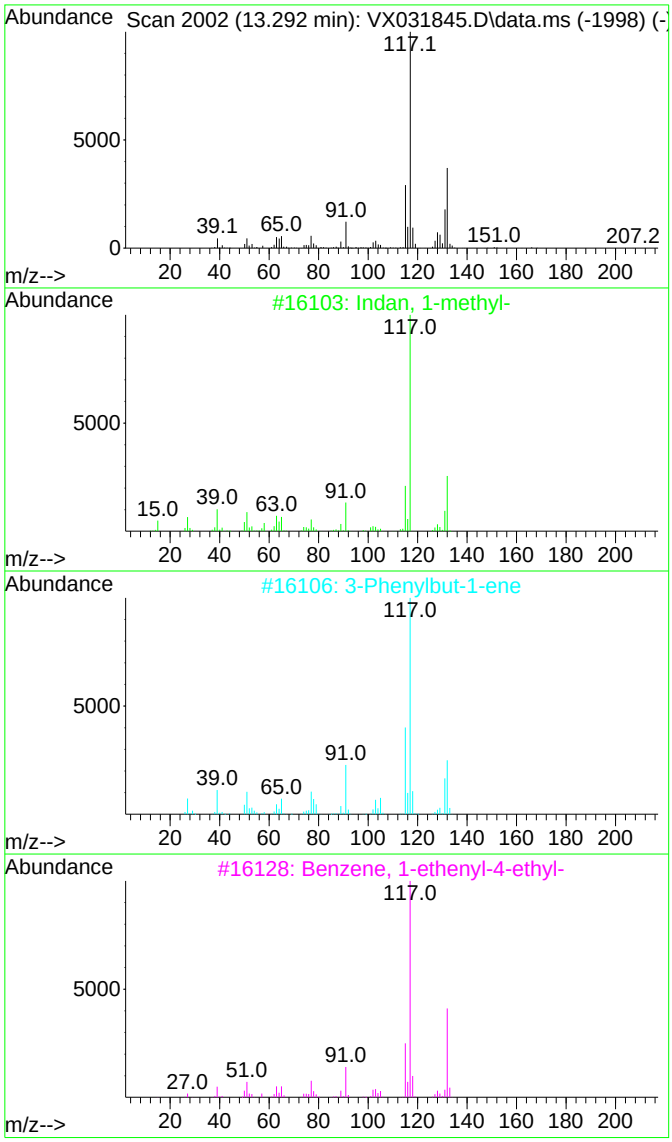
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100322W.M
Quant Title : SW846 8260

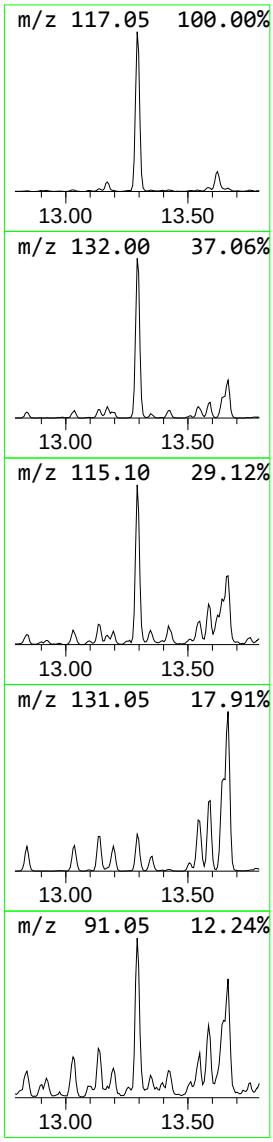
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Indan, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	48.07 ug/l	403162	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	90
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
3	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
5	1-Phenyl-1-butene	132	C10H12	000824-90-8	86





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100322\
Data File : VX031845.D
Acq On : 03 Oct 2022 19:48
Operator : JC/MD
Sample : N4860-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2R

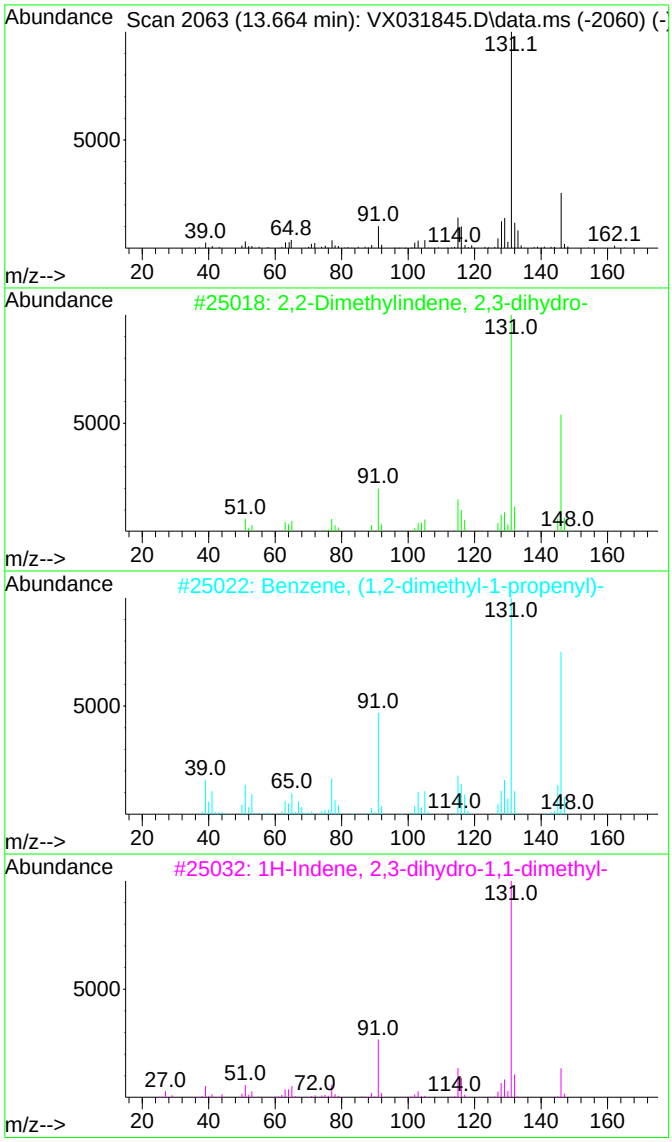
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100322W.M
Quant Title : SW846 8260

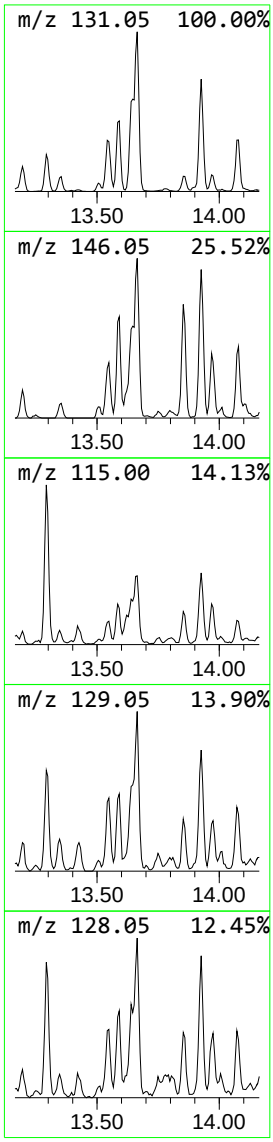
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 2,2-Dimethylindene, 2,3-dih... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.664	33.10 ug/l	277605	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	91
2	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	91
3	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	91
4	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	90
5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	90





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Data File : VX031845.D
Acq On : 03 Oct 2022 19:48
Operator : JC/MD
Sample : N4860-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2R

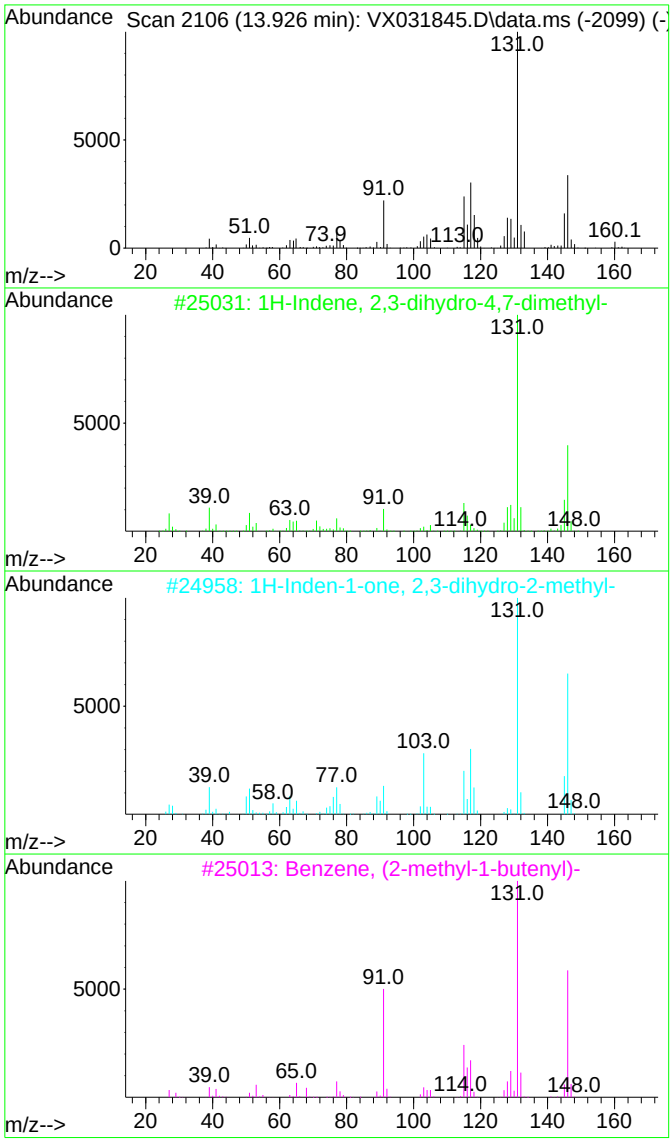
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100322W.M
Quant Title : SW846 8260

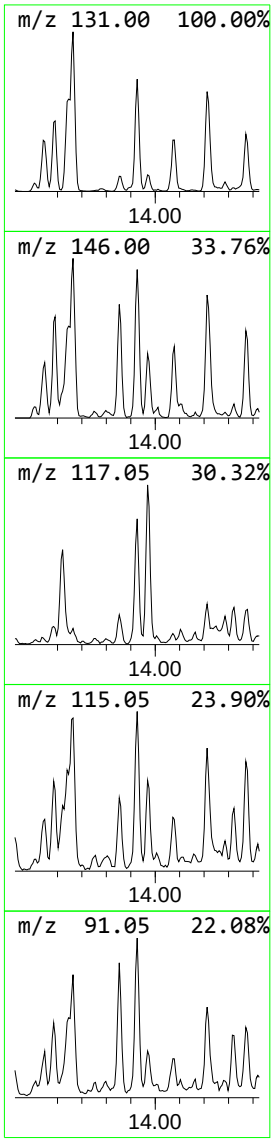
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.926	35.58 ug/l	298432	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
2	1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	90
3	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	90
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	87
5	Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	81





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100322\
Data File : VX031845.D
Acq On : 03 Oct 2022 19:48
Operator : JC/MD
Sample : N4860-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2R

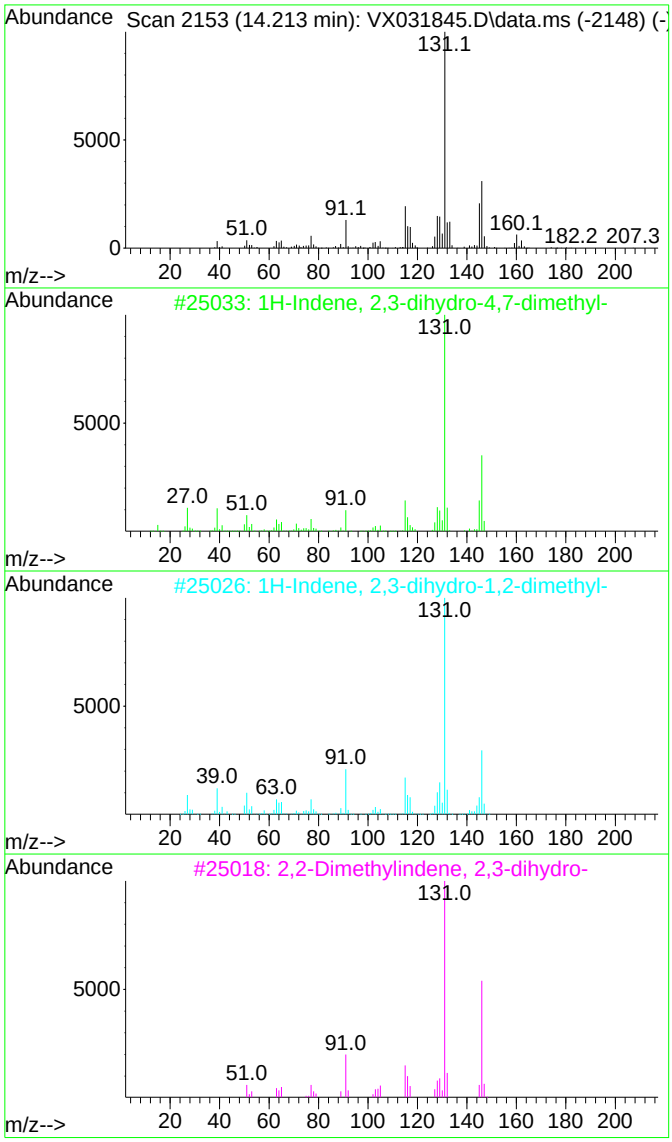
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100322W.M
Quant Title : SW846 8260

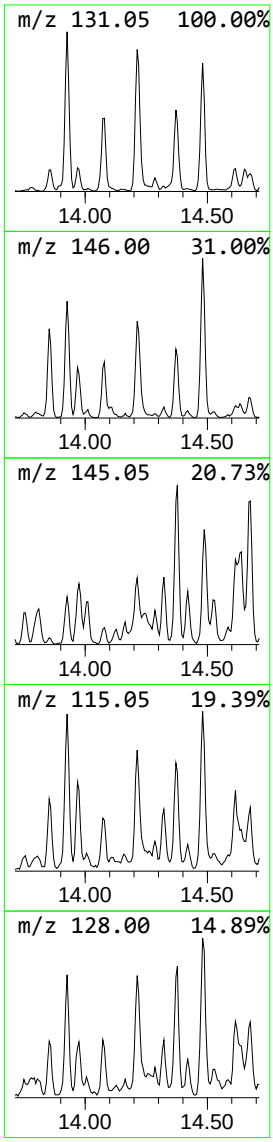
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.213	35.46 ug/l	297420	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91
2	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91
3	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	87
4	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	87
5	1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14	001685-82-1	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100322\
Data File : VX031845.D
Acq On : 03 Oct 2022 19:48
Operator : JC/MD
Sample : N4860-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2R

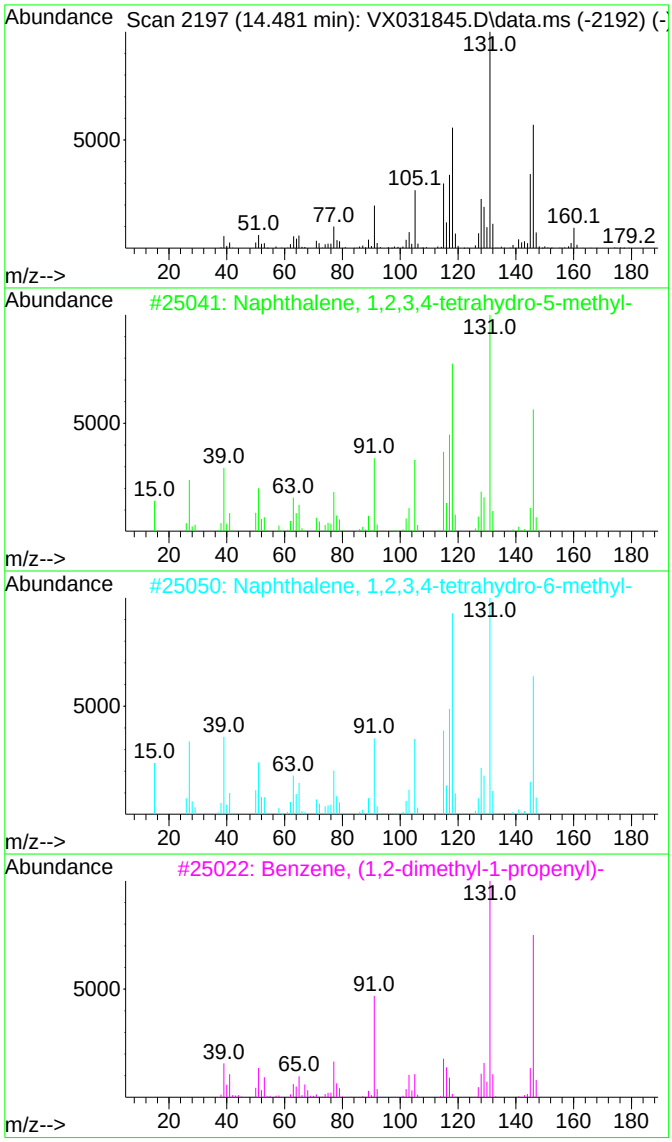
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100322W.M
Quant Title : SW846 8260

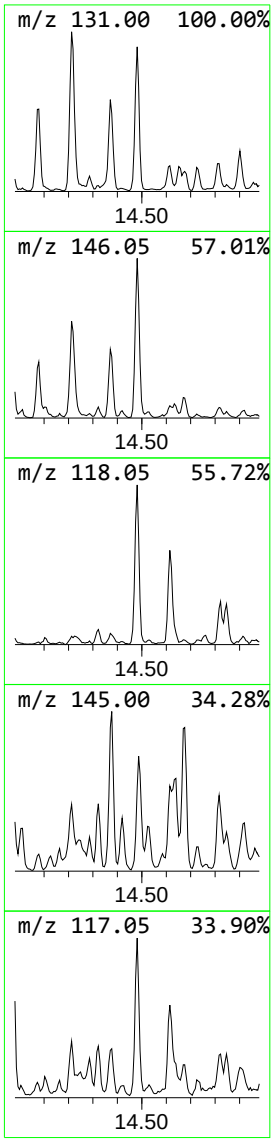
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.481	42.72 ug/l	358287	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	93
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	89
3	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	60
4	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	58
5	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	58





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100322\
Data File : VX031845.D
Acq On : 03 Oct 2022 19:48
Operator : JC/MD
Sample : N4860-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2R

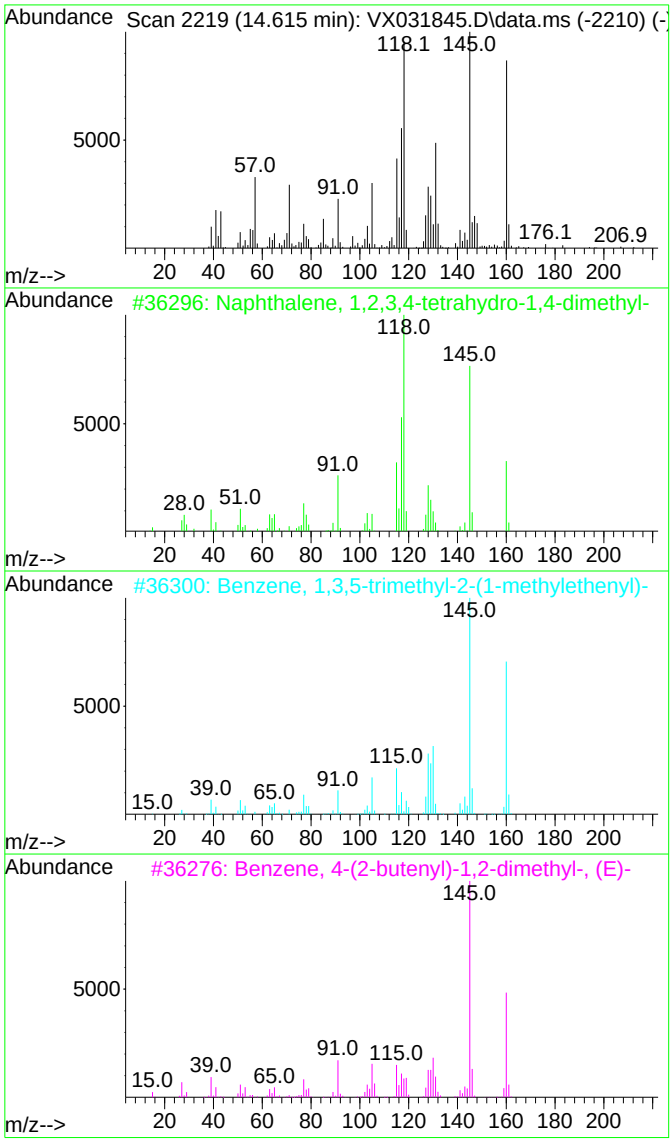
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100322W.M
Quant Title : SW846 8260

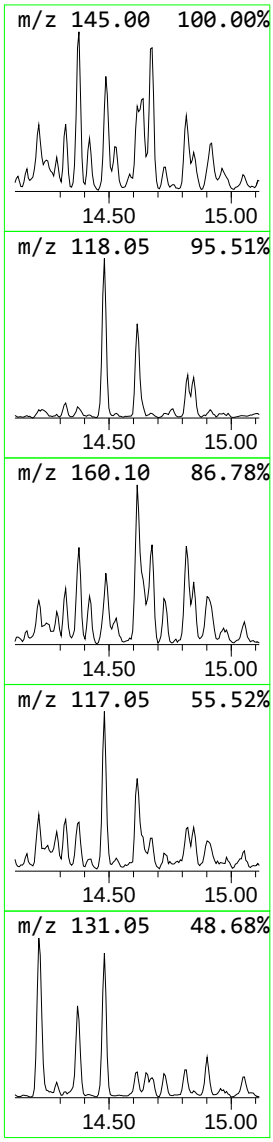
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1,3,5-trimethyl-2-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.615	37.88 ug/l	317735	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	95
2	Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	70
3	Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	64
4	Benzene, 1-(1-methylethenyl)-3-(...	160	C12H16	001129-29-9	58
5	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	58





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100322\
Data File : VX031845.D
Acq On : 03 Oct 2022 19:48
Operator : JC/MD
Sample : N4860-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2R

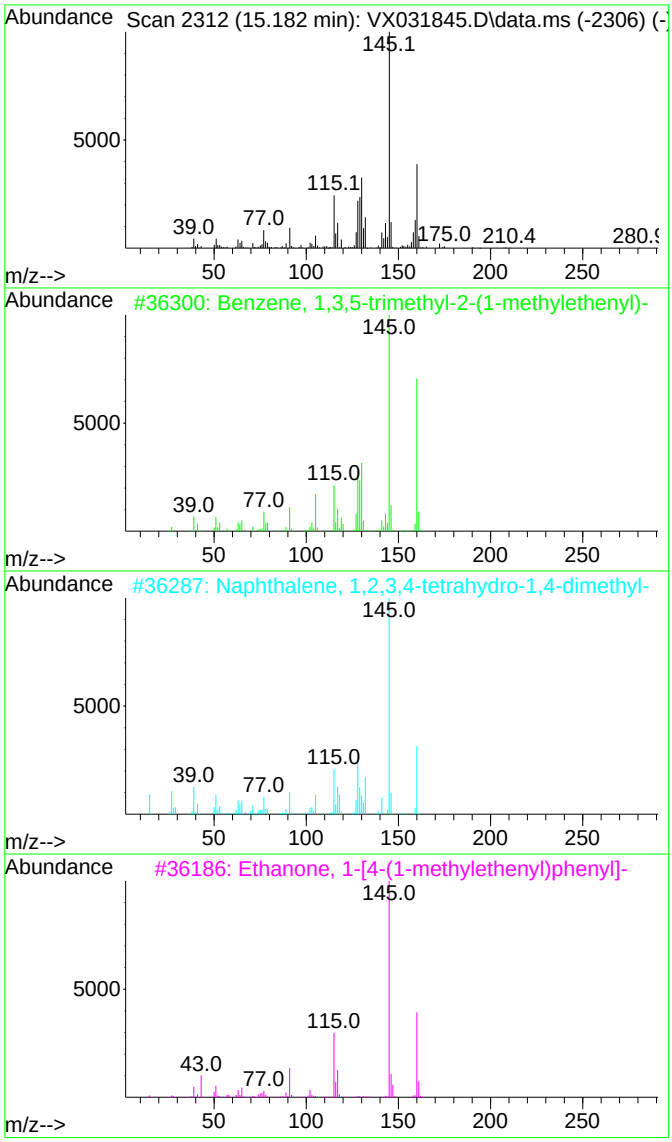
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100322W.M
Quant Title : SW846 8260

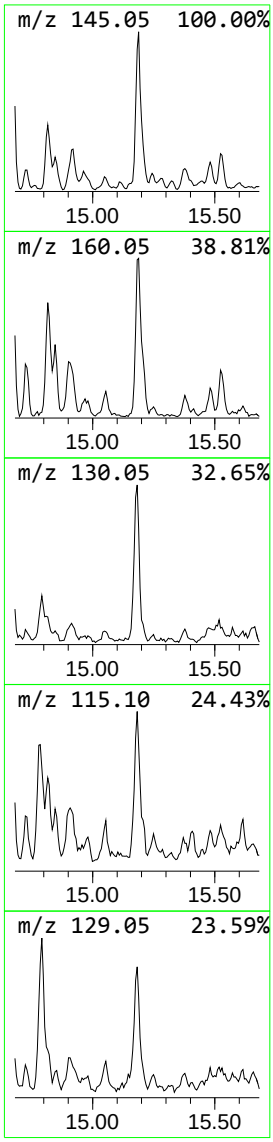
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.182	29.06 ug/l	243747	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	95
2	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	94
3	Ethanone, 1-[4-(1-methylethenyl)...	160	C11H12O	005359-04-6	90
4	Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	87
5	1H-Indene, 2,3-dihydro-1,1,3-tri...	160	C12H16	002613-76-5	68





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100322\
Data File : VX031845.D
Acq On : 03 Oct 2022 19:48
Operator : JC/MD
Sample : N4860-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100322W.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	1.282	53.2	ug/l	234239	1	5.556	220287	50.0
Indane	12.243	75.2	ug/l	630725	4	12.024	419344	50.0
Benzene, 1-ethe...	12.701	36.0	ug/l	301965	4	12.024	419344	50.0
Indan, 1-methyl-	13.292	48.1	ug/l	403162	4	12.024	419344	50.0
2,2-Dimethylind...	13.664	33.1	ug/l	277605	4	12.024	419344	50.0
1H-Indene, 2,3-...	13.926	35.6	ug/l	298432	4	12.024	419344	50.0
1H-Indene, 2,3-...	14.213	35.5	ug/l	297420	4	12.024	419344	50.0
Naphthalene, 1,...	14.481	42.7	ug/l	358287	4	12.024	419344	50.0
Benzene, 1,3,5-...	14.615	37.9	ug/l	317735	4	12.024	419344	50.0
Naphthalene, 1,...	15.182	29.1	ug/l	243747	4	12.024	419344	50.0