LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062921\
Data File : VX023004.D
Acq On : 29 Jun 2021 15:32
Operator : JC/MD
Sample : M2875-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-5

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\82X062321W.M
Title : SW846 8260

Signal : TIC: VX023004.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.117	322	333	346	rVB3	74599	193543	5.98%	1.152%
2	4.306	520	528	540	rVB	51852	148118	4.57%	0.882%
3	5.397	694	707	714	rBV	46168	139992	4.32%	0.833%
4	5.483	714	721	727	rBV4	36466	113345	3.50%	0.675%
5	5.562	728	734	741	rVB2	53006	134784	4.16%	0.802%
6	5.836	767	779	791	rBV3	41280	128302	3.96%	0.764%
7	5.964	791	800	806	rBV	62820	161405	4.98%	0.961%
8	6.050	806	814	824	rVB	322822	836276	25.83%	4.978%
9	6.178	824	835	836	rBV3	28489	70367	2.17%	0.419%
10	6.263	844	849	855	rVB2	37493	83425	2.58%	0.497%
11	6.342	855	862	878	rVB5	76463	266283	8.22%	1.585%
12	6.611	897	906	921	rVB4	13230	35066	1.08%	0.209%
13	6.769	921	932	939	rBV	133335	335325	10.36%	1.996%
14	6.855	940	946	957	rVB4	46156	142658	4.41%	0.849%
15	7.068	975	981	991	rVV3	15085	36568	1.13%	0.218%
16	7.178	991	999	1008	rVB2	24305	53777	1.66%	0.320%
17	7.385	1020	1033	1045	rBV5	158807	438365	13.54%	2.609%
18	7.665	1072	1079	1088	rVB	39767	77504	2.39%	0.461%
19	7.860	1104	1111	1112	rBV	32003	49570	1.53%	0.295%
20	7.885	1113	1115	1122	rVB3	36752	59493	1.84%	0.354%
21	7.988	1122	1132	1144	rVB2	36429	97139	3.00%	0.578%
22	8.110	1144	1152	1155	rBV2	54975	111537	3.44%	0.664%
23	8.147	1155	1158	1162	rVB	39561	59687	1.84%	0.355%
24	8.312	1181	1185	1194	rVB7	15918	33123	1.02%	0.197%
25	8.446	1201	1207	1211	rBV5	15409	33916	1.05%	0.202%
26	8.507	1211	1217	1228	rVB3	85271	190936	5.90%	1.136%
27	8.604	1228	1233	1236	rBV3	43044	78474	2.42%	0.467%
28	8.653	1236	1241	1249	rVB	274773	446673	13.79%	2.659%
29	8.927	1281	1286	1290	rBV	22786	41554	1.28%	0.247%
30	8.976	1290	1294	1301	rVB3	32173	59976	1.85%	0.357%
31	9.104	1309	1315	1320	rVB	39586	65576	2.03%	0.390%
32	9.165	1320	1325	1336	rVB	86511	145816	4.50%	0.868%
33	9.513	1377	1382	1387	rVB4	32240	65818	2.03%	0.392%
34	9.567	1387	1391	1395	rVV2	30188	42380	1.31%	0.252%
35	9.763	1417	1423	1430	rBV4	26658	48345	1.49%	0.288%
36	10.055	1461	1471	1476	rBV	270753	377039	11.64%	2.244%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062921\
Data File : VX023004.D
Acq On : 29 Jun 2021 15:32
Operator : JC/MD
Sample : M2875-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-5

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\82X062321W.M
Title : SW846 8260

37	10.134	1481	1484	1491	rVV4	30292	66561	2.06%	0.396%
38	10.305	1505	1512	1518	rVB	202399	283984	8.77%	1.690%
39	10.647	1563	1568	1573	rVB	27720	34512	1.07%	0.205%
40	10.964	1615	1620	1629	rBV	237135	297030	9.17%	1.768%
41	11.085	1635	1640	1645	rVB	214715	265600	8.20%	1.581%
42	11.305	1672	1676	1683	rVB	244492	315829	9.75%	1.880%
43	11.402	1685	1692	1696	rVV	67413	94814	2.93%	0.564%
44	11.457	1696	1701	1707	rVB2	18286	33534	1.04%	0.200%
45	11.616	1722	1727	1735	rVB	343324	419064	12.94%	2.494%
46	11.756	1746	1750	1755	rVB	27363	33532	1.04%	0.200%
47	11.866	1763	1768	1770	rBV	45170	56771	1.75%	0.338%
48	11.890	1770	1772	1780	rVB	58549	72537	2.24%	0.432%
49	12.024	1788	1794	1799	rBV	270788	332767	10.28%	1.981%
50	12.091	1801	1805	1811	rVB	67168	83163	2.57%	0.495%
51	12.244	1825	1830	1837	rBV2	2454155	3238152	100.00%	19.274%
52	12.317	1837	1842	1853	rVV2	167550	258504	7.98%	1.539%
53	12.408	1853	1857	1862	rVV2	68023	93885	2.90%	0.559%
54	12.469	1862	1867	1873	rVB3	110484	159895	4.94%	0.952%
55	12.616	1886	1891	1894	rBV	607093	737285	22.77%	4.388%
56	12.646	1894	1896	1901	rVV	188313	218273	6.74%	1.299%
57	12.701	1901	1905	1913	rVB2	801501	1055782	32.60%	6.284%
58	12.847	1924	1929	1933	rBV2	29540	48137	1.49%	0.287%
59	12.927	1937	1942	1946	rVV	397517	481331	14.86%	2.865%
60	12.969	1946	1949	1954	rVV3	25939	40173	1.24%	0.239%
61	13.030	1954	1959	1964	rVB3	39760	61670	1.90%	0.367%
62	13.170	1978	1982	1991	rVB	382735	459624	14.19%	2.736%
63	13.292	1991	2002	2008	rVV2	594236	845892	26.12%	5.035%
64	13.347	2008	2011	2013	rVV2	43256	55913	1.73%	0.333%
65	13.372	2013	2015	2020	rVV	27327	35744	1.10%	0.213%
66	13.427	2020	2024	2032	rVB	73453	96183	2.97%	0.572%
67	13.506	2032	2037	2040	rBV	25347	33331	1.03%	0.198%
68	13.548	2040	2044	2047	rVV	94245	131866	4.07%	0.785%
69	13.585	2047	2050	2054	rVV2	65249	92070	2.84%	0.548%
70	13.646	2054	2060	2061	rVV2	53336	86121	2.66%	0.513%
71	13.664	2061	2063	2072	rVB2	81467	117352	3.62%	0.698%
72	13.780	2072	2082	2090	rBV	175515	246304	7.61%	1.466%
73	13.926	2100	2106	2110	rBV3	29776	41845	1.29%	0.249%
74	14.079	2126	2131	2138	rBV	38410	51530	1.59%	0.307%
75	14.219	2149	2154	2160	rBV3	30555	59159	1.83%	0.352%
76	14.640	2218	2223	2227	rVB	67988	82063	2.53%	0.488%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062921\
Data File : VX023004.D
Acq On : 29 Jun 2021 15:32
Operator : JC/MD
Sample : M2875-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-5

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\82X062321W.M
Title : SW846 8260

77	14.780	2240	2246	2252	rBV	98663	129921	4.01%	0.773%
78	16.328	2493	2500	2511	rBV	26801	51123	1.58%	0.304%

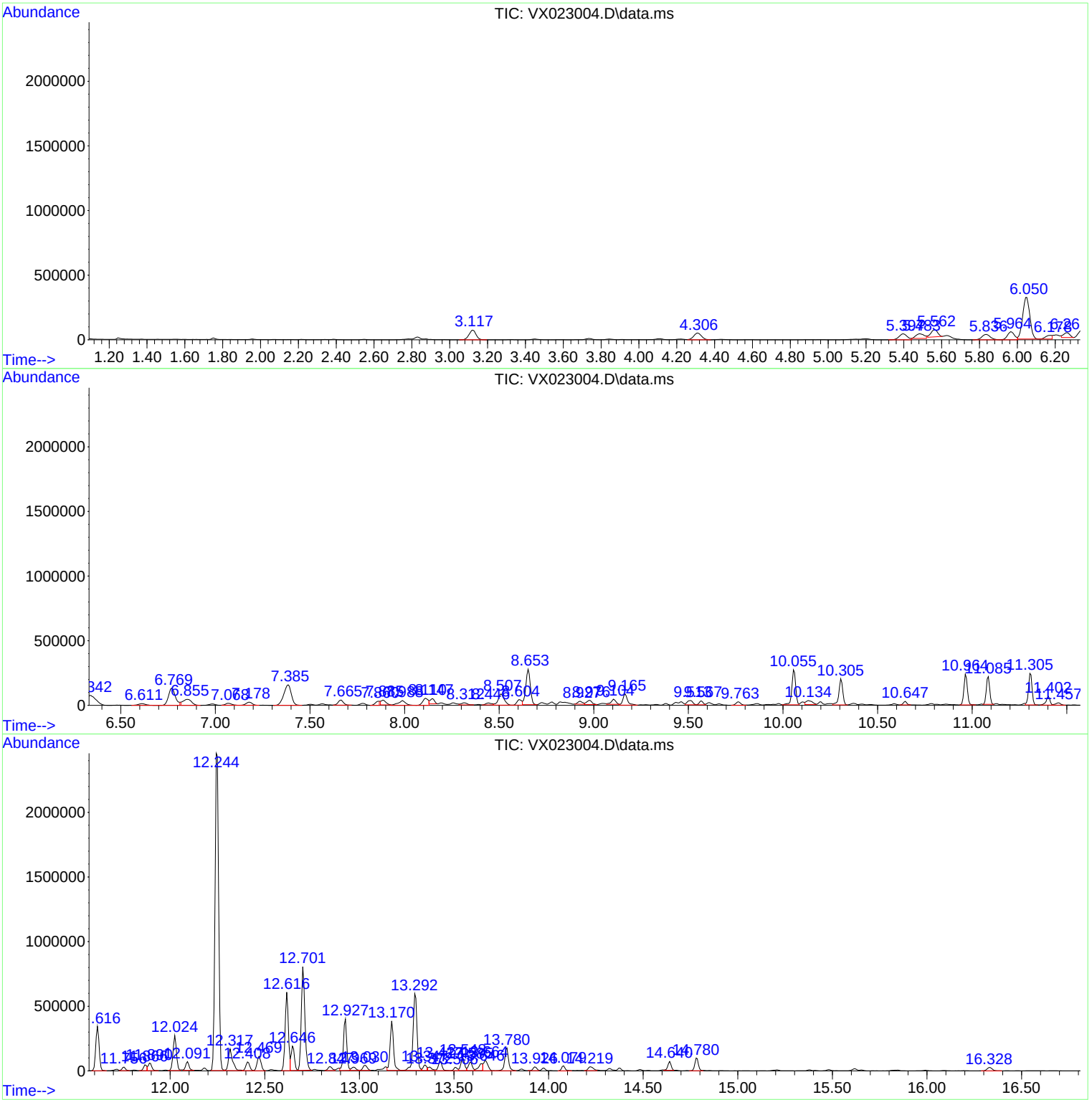
Sum of corrected areas: 16800981

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062921\
Data File : VX023004.D
Acq On : 29 Jun 2021 15:32
Operator : JC/MD
Sample : M2875-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-5

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062921\
Data File : VX023004.D
Acq On : 29 Jun 2021 15:32
Operator : JC/MD
Sample : M2875-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-5

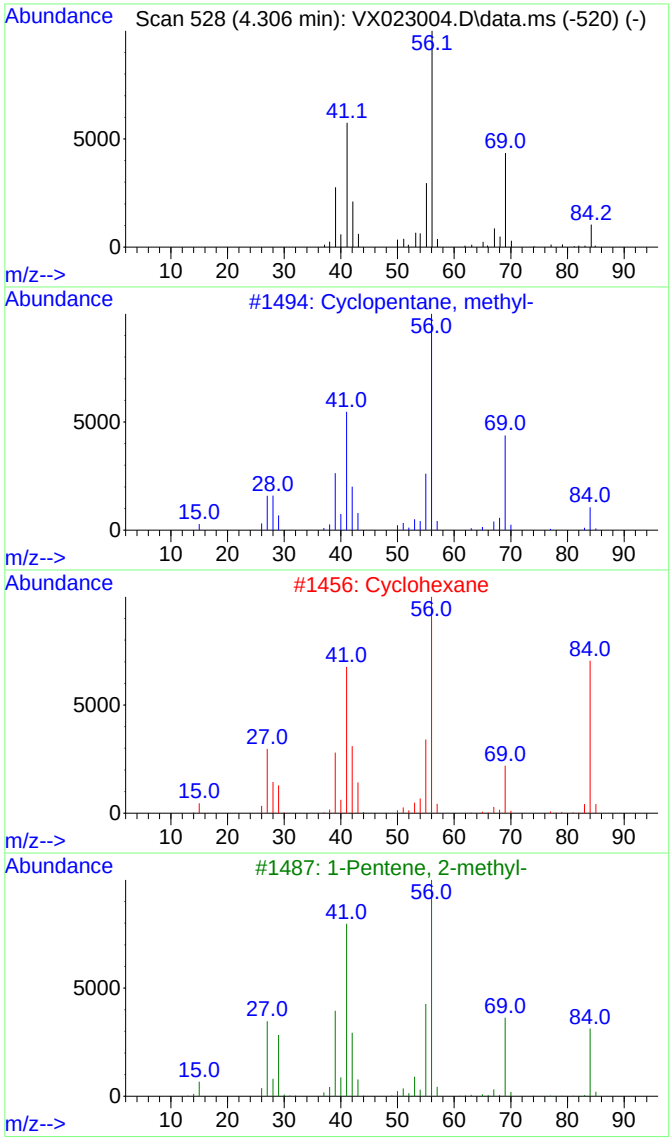
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
Quant Title : SW846 8260

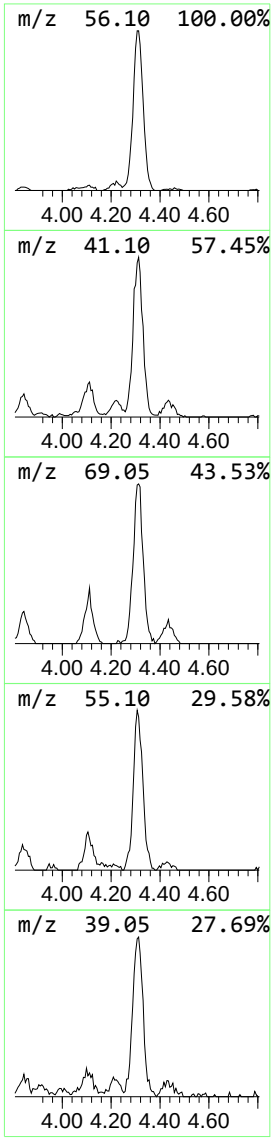
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Cyclopentane, methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.306	54.95 ug/l	148118	Pentafluorobenzene	5.562

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2	Cyclohexane	84	C6H12	000110-82-7	90
3	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	83
4	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
5	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062921\
Data File : VX023004.D
Acq On : 29 Jun 2021 15:32
Operator : JC/MD
Sample : M2875-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-5

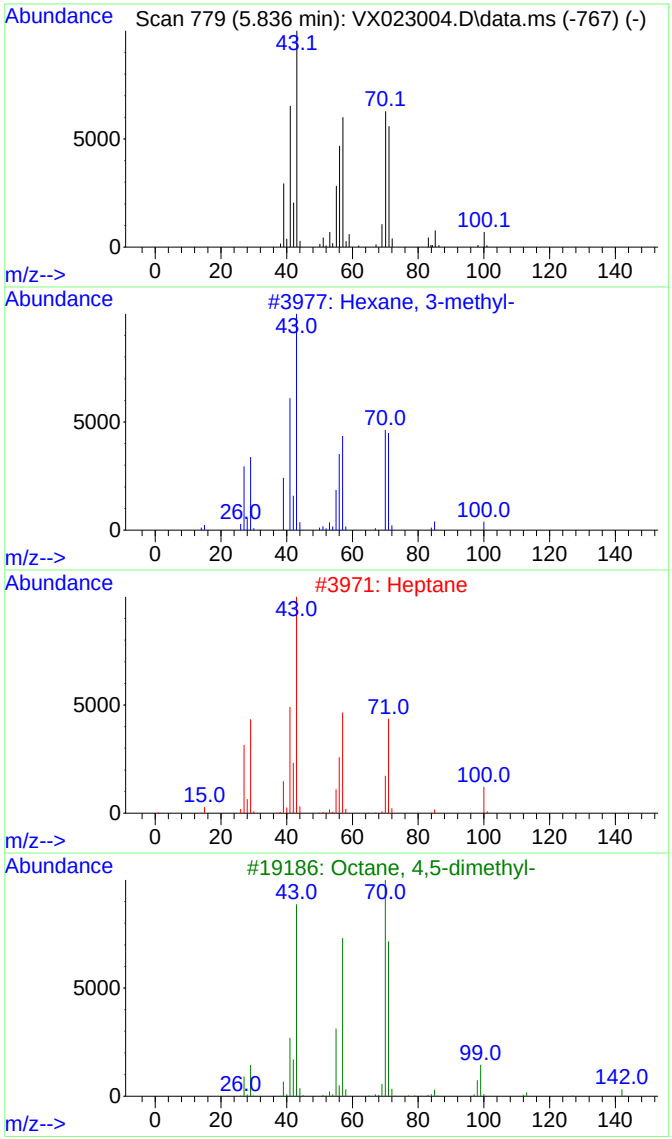
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
Quant Title : SW846 8260

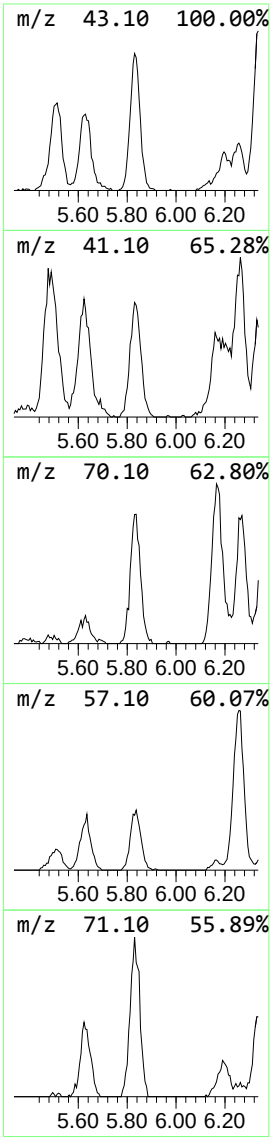
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.836	47.60 ug/l	128302	Pentafluorobenzene	5.562

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 3-methyl-	100	C7H16	000589-34-4	93
2	Heptane	100	C7H16	000142-82-5	72
3	Octane, 4,5-dimethyl-	142	C10H22	015869-96-2	53
4	Pyrrolidine	71	C4H9N	000123-75-1	47
5	Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	40





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062921\
Data File : VX023004.D
Acq On : 29 Jun 2021 15:32
Operator : JC/MD
Sample : M2875-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-5

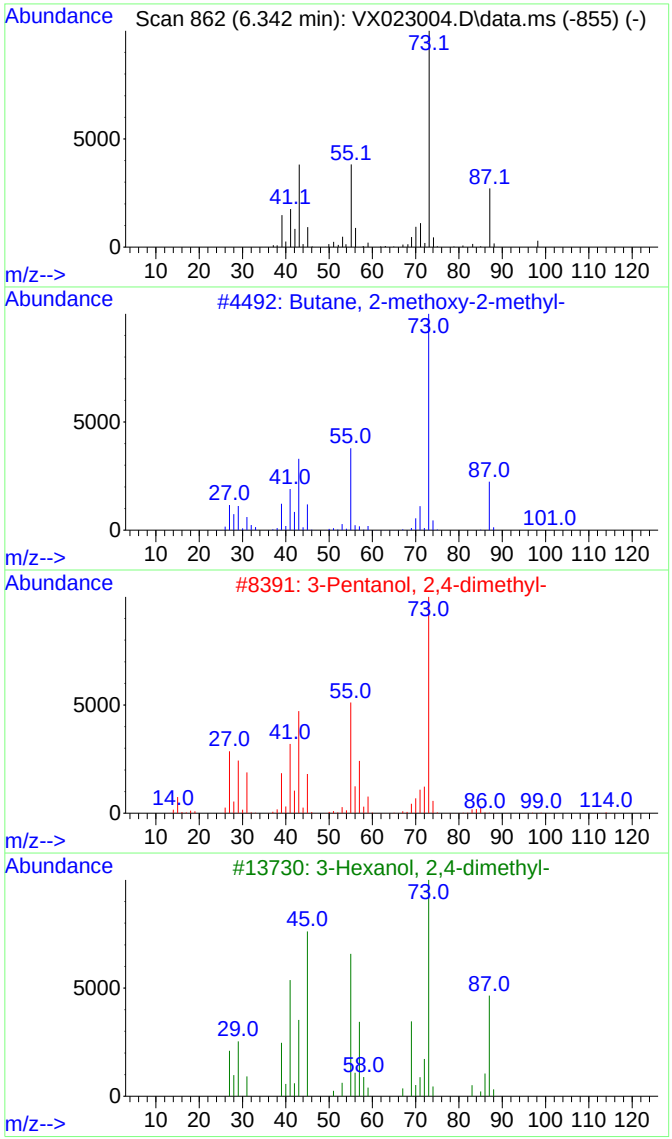
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
Quant Title : SW846 8260

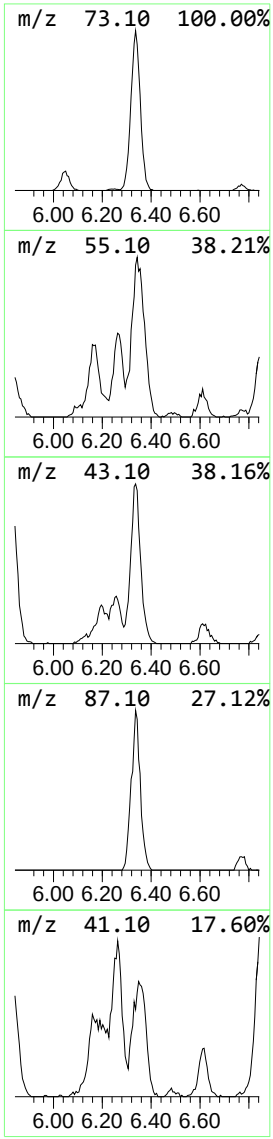
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Butane, 2-methoxy-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.342	39.71 ug/l	266283	1,4-Difluorobenzene	6.769

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	59
2	3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	50
3	3-Hexanol, 2,4-dimethyl-	130	C8H18O	013432-25-2	28
4	2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	23
5	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	10





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062921\
Data File : VX023004.D
Acq On : 29 Jun 2021 15:32
Operator : JC/MD
Sample : M2875-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-5

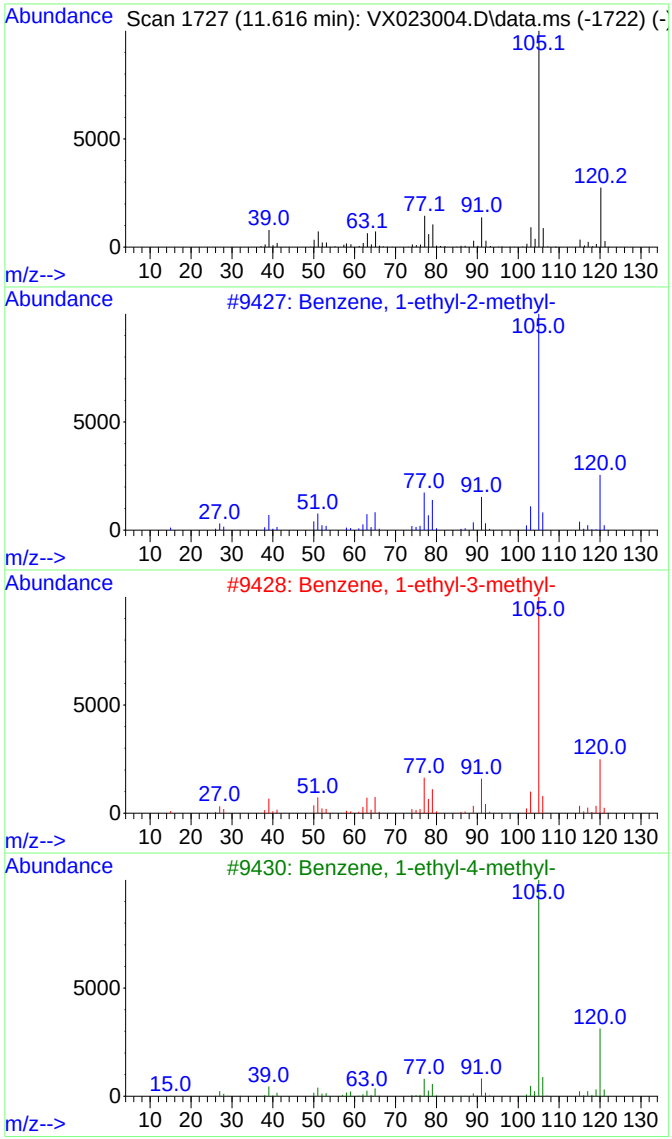
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
Quant Title : SW846 8260

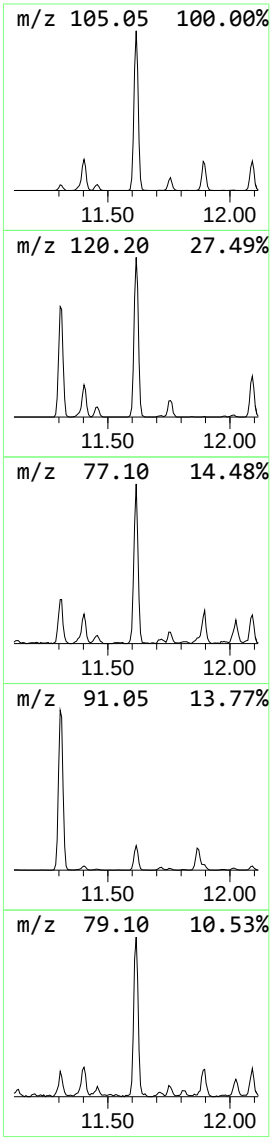
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.616	62.97 ug/l	419064	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	93
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	93
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4	Mesitylene	120	C9H12	000108-67-8	91
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062921\
Data File : VX023004.D
Acq On : 29 Jun 2021 15:32
Operator : JC/MD
Sample : M2875-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-5

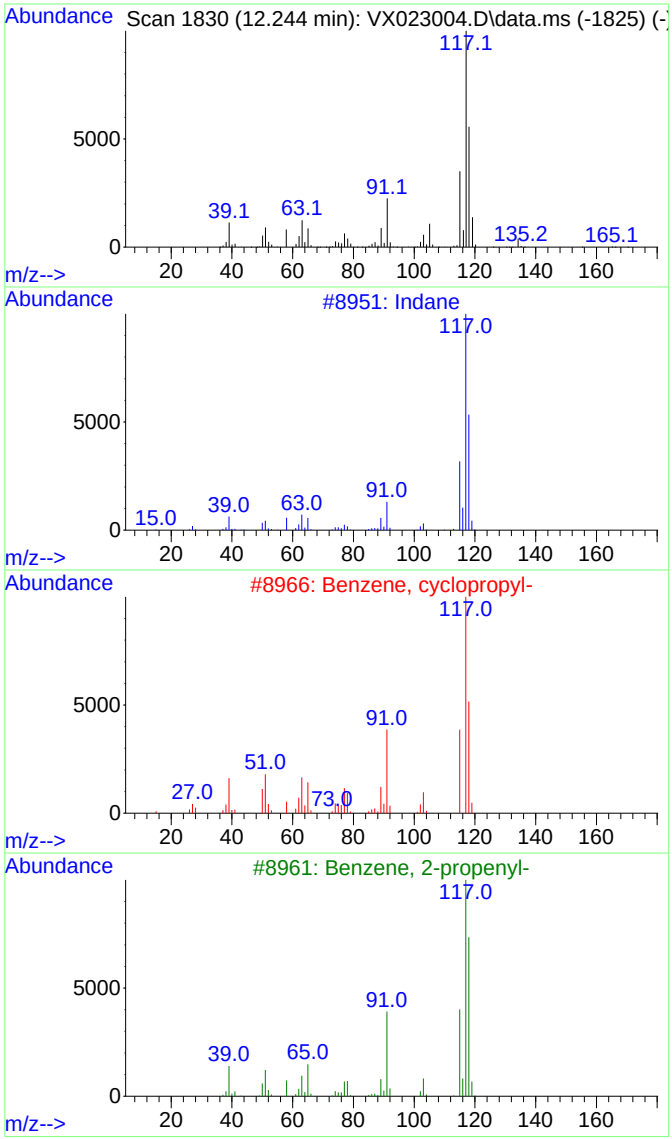
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
Quant Title : SW846 8260

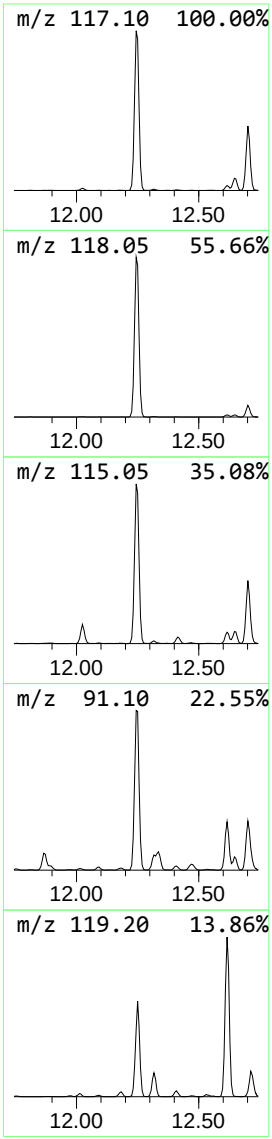
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	486.55 ug/l	3238150	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	76
4	Deltacyclene	118	C9H10	007785-10-6	58
5	(Z)-1-Phenylpropene	118	C9H10	000766-90-5	55





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062921\
Data File : VX023004.D
Acq On : 29 Jun 2021 15:32
Operator : JC/MD
Sample : M2875-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-5

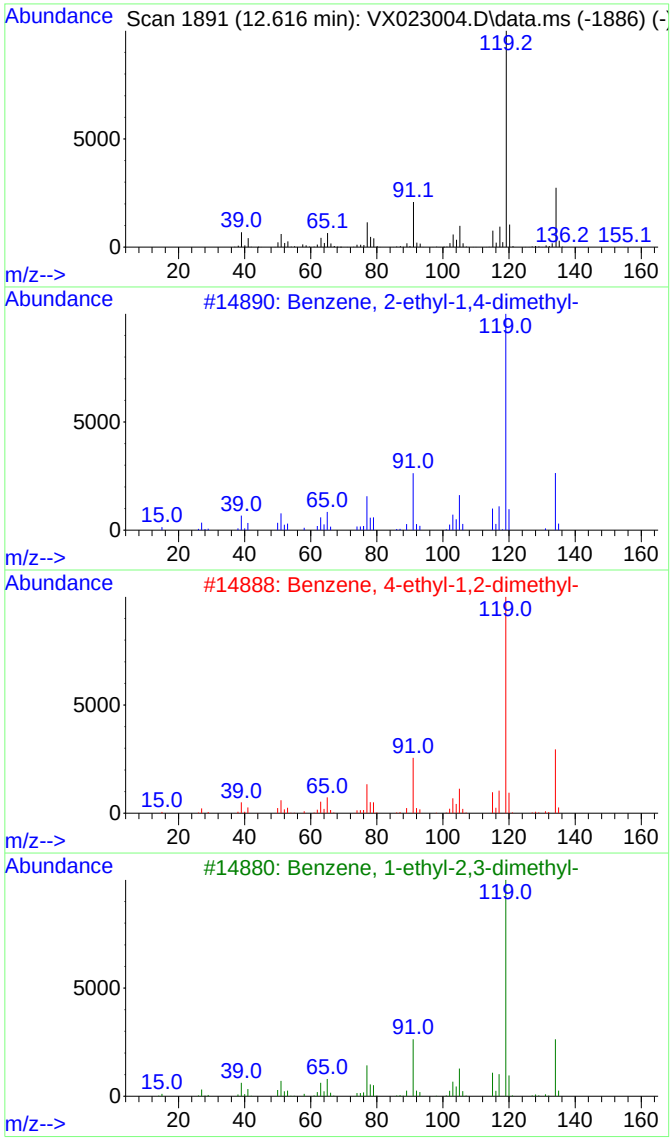
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
Quant Title : SW846 8260

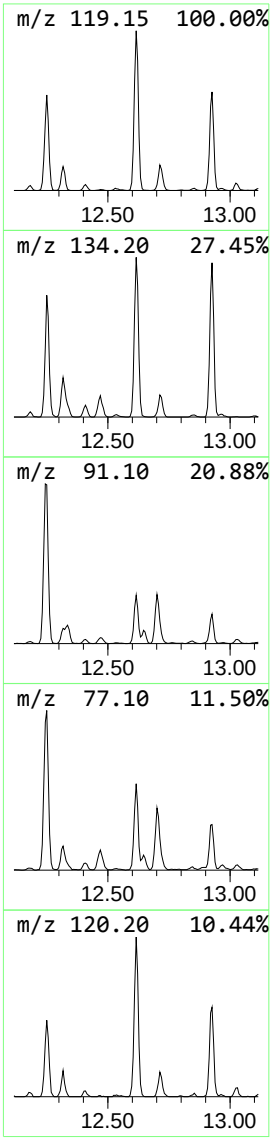
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.616	110.78 ug/l	737285	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
5	o-Cymene	134	C10H14	000527-84-4	95





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062921\
Data File : VX023004.D
Acq On : 29 Jun 2021 15:32
Operator : JC/MD
Sample : M2875-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-5

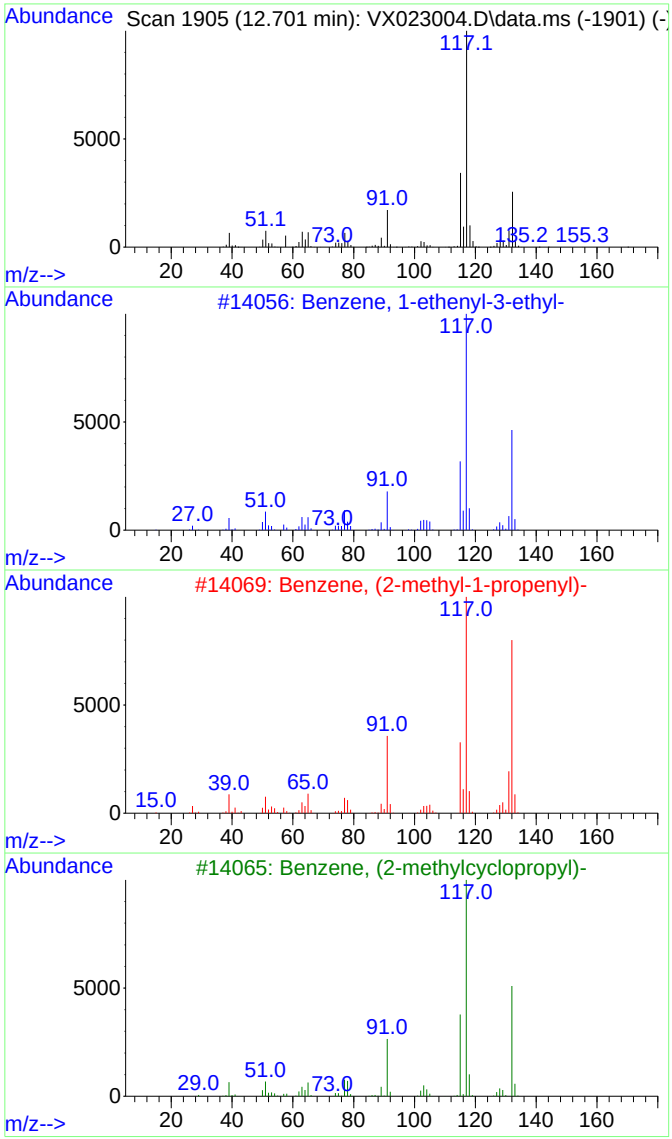
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
Quant Title : SW846 8260

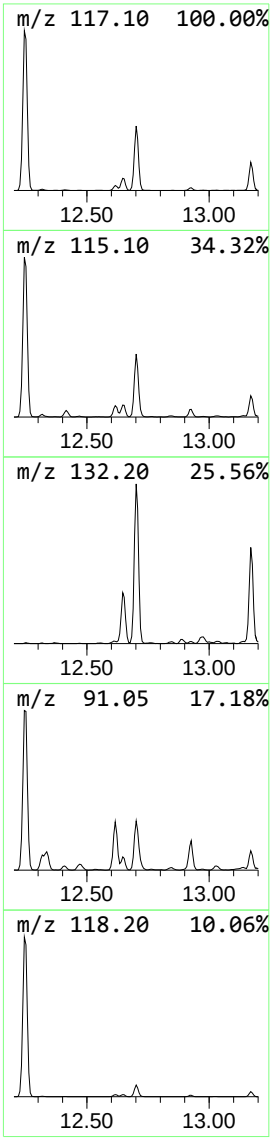
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	158.64 ug/l	1055780	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	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90
3	Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	87
4	Indan, 1-methyl-	132	C10H12	000767-58-8	87
5	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	83





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062921\
Data File : VX023004.D
Acq On : 29 Jun 2021 15:32
Operator : JC/MD
Sample : M2875-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-5

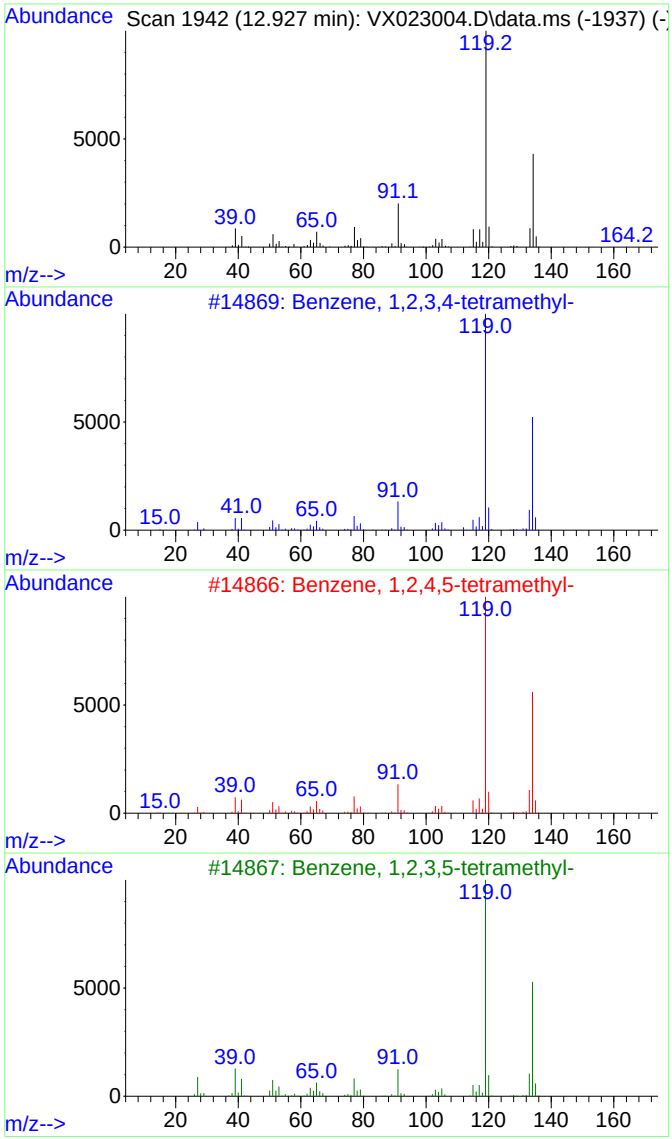
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
Quant Title : SW846 8260

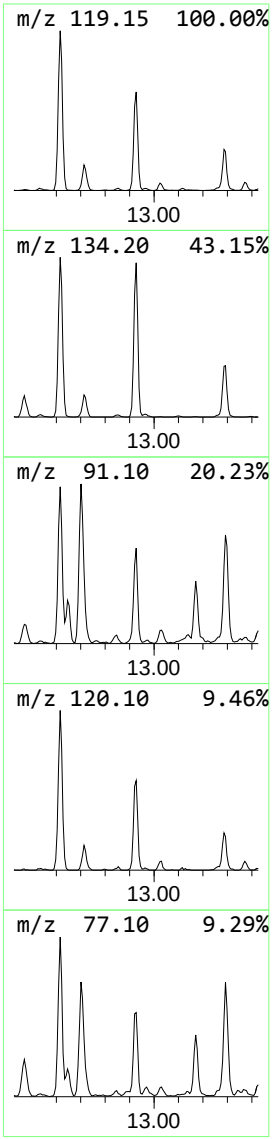
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.927	72.32 ug/l	481331	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062921\
Data File : VX023004.D
Acq On : 29 Jun 2021 15:32
Operator : JC/MD
Sample : M2875-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-5

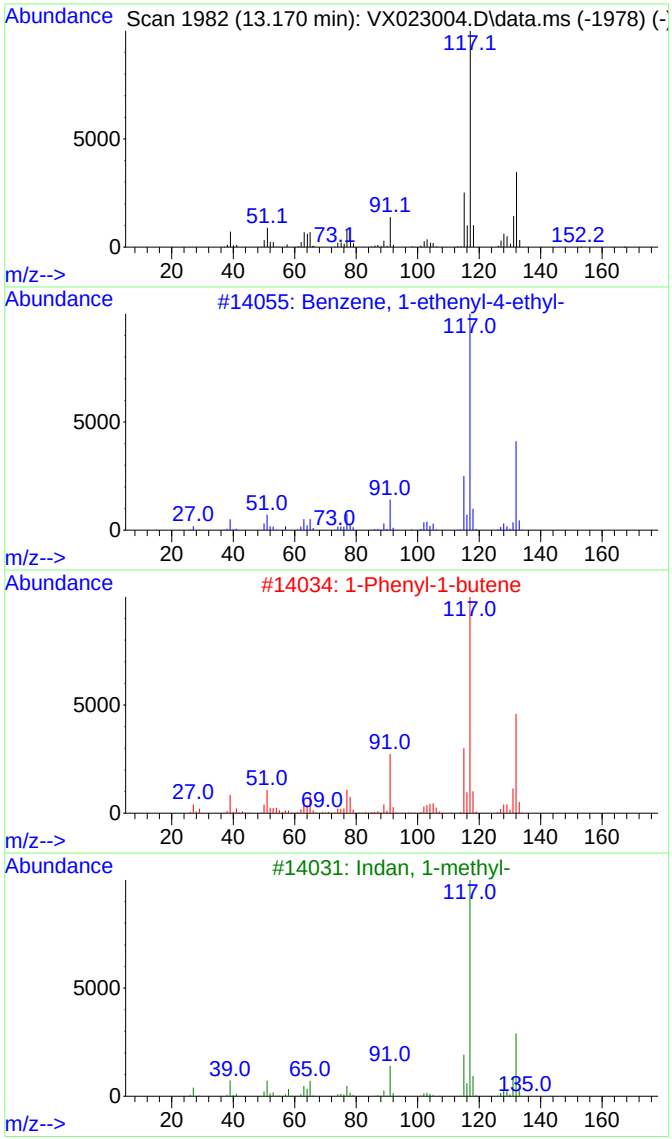
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
Quant Title : SW846 8260

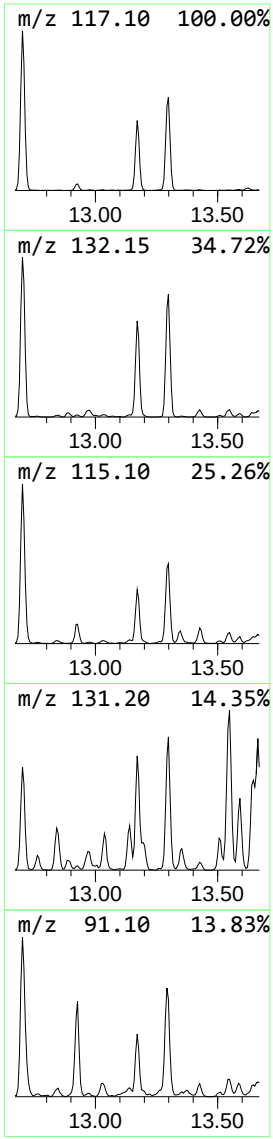
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.171	69.06 ug/l	459624	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	94
2	1-Phenyl-1-butene	132	C10H12	000824-90-8	93
3	Indan, 1-methyl-	132	C10H12	000767-58-8	91
4	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
5	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062921\
Data File : VX023004.D
Acq On : 29 Jun 2021 15:32
Operator : JC/MD
Sample : M2875-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-5

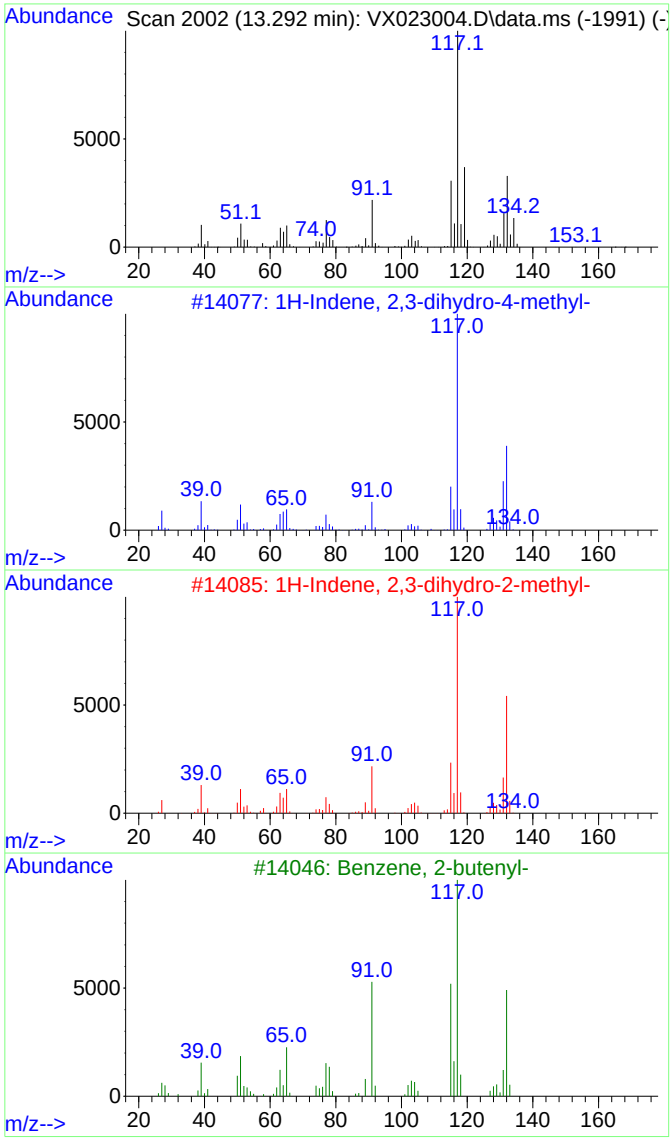
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
Quant Title : SW846 8260

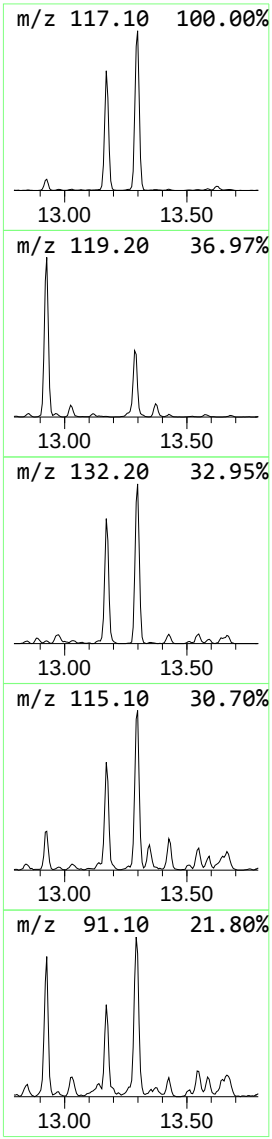
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	127.10 ug/l	845892	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	93
2	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	91
3	Benzene, 2-butenyl-	132	C10H12	001560-06-1	90
4	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	90
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062921\
Data File : VX023004.D
Acq On : 29 Jun 2021 15:32
Operator : JC/MD
Sample : M2875-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-5

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Cyclopentane, m...	4.306	55.0	ug/l	148118	1	5.562	134784	50.0
Hexane, 3-methyl-	5.836	47.6	ug/l	128302	1	5.562	134784	50.0
Butane, 2-metho...	6.342	39.7	ug/l	266283	2	6.769	335325	50.0
Benzene, 1-ethy...	11.616	63.0	ug/l	419064	4	12.024	332767	50.0
Indane	12.244	486.6	ug/l	3238150	4	12.024	332767	50.0
Benzene, 2-ethy...	12.616	110.8	ug/l	737285	4	12.024	332767	50.0
Benzene, 1-ethe...	12.701	158.6	ug/l	1055780	4	12.024	332767	50.0
Benzene, 1,2,3,...	12.927	72.3	ug/l	481331	4	12.024	332767	50.0
Benzene, 1-ethe...	13.171	69.1	ug/l	459624	4	12.024	332767	50.0
1H-Indene, 2,3-...	13.292	127.1	ug/l	845892	4	12.024	332767	50.0