LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111522\
Data File : VX032826.D
Acq On : 15 Nov 2022 16:35
Operator : JC/MD
Sample : N5614-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

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\82X102622W.M
Title : SW846 8260

Signal : TIC: VX032826.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	41	47	53	rBV	66556	74109	1.29%	0.219%
2	1.752	104	109	120	rVB	156454	208381	3.63%	0.615%
3	2.782	259	278	282	rBV2	60602	151220	2.63%	0.447%
4	2.825	282	285	290	rVV	72512	148554	2.59%	0.439%
5	2.867	290	292	304	rVB3	31766	64452	1.12%	0.190%
6	3.105	322	331	342	rVB2	141029	323389	5.63%	0.955%
7	3.751	425	437	445	rBV3	45371	155376	2.71%	0.459%
8	4.208	503	512	520	rVV2	29028	85866	1.50%	0.254%
9	4.306	520	528	540	rVB2	83631	236614	4.12%	0.699%
10	5.190	650	673	690	rBV4	27716	120180	2.09%	0.355%
11	5.385	694	705	713	rVV2	67443	199762	3.48%	0.590%
12	5.489	713	722	725	rVV5	41912	142314	2.48%	0.420%
13	5.550	726	732	738	rVV	106637	319360	5.56%	0.943%
14	5.623	739	744	763	rVB3	67129	213227	3.71%	0.630%
15	5.830	765	778	790	rBV3	53850	165195	2.88%	0.488%
16	5.958	790	799	804	rVV	66474	177493	3.09%	0.524%
17	6.037	804	812	825	rVV	819547	2191789	38.18%	6.472%
18	6.159	825	832	839	rVV3	60931	207204	3.61%	0.612%
19	6.257	840	848	857	rVV2	139268	438138	7.63%	1.294%
20	6.360	857	865	877	rVB3	49720	149587	2.61%	0.442%
21	6.763	920	931	941	rBV	160579	448213	7.81%	1.324%
22	6.854	942	946	957	rVB3	46431	111160	1.94%	0.328%
23	7.171	990	998	1008	rBV2	76451	168418	2.93%	0.497%
24	7.360	1019	1029	1044	rVB4	74674	208872	3.64%	0.617%
25	7.885	1104	1115	1122	rBV6	33804	123587	2.15%	0.365%
26	7.988	1122	1132	1142	rVB2	90388	225339	3.93%	0.665%
27	8.104	1142	1151	1156	rBV	167973	377310	6.57%	1.114%
28	8.189	1162	1165	1178	rVB5	33483	74135	1.29%	0.219%
29	8.311	1178	1185	1190	rBV2	40291	79772	1.39%	0.236%
30	8.506	1209	1217	1226	rVB3	117718	229833	4.00%	0.679%
31	8.604	1226	1233	1235	rBV2	59477	103559	1.80%	0.306%
32	8.647	1235	1240	1248	rVB	365296	601477	10.48%	1.776%
33	9.104	1307	1315	1320	rBV	60367	117297	2.04%	0.346%
34	9.159	1320	1324	1335	rVB	187158	306306	5.34%	0.904%
35	9.433	1364	1369	1370	rBV	48662	69362	1.21%	0.205%
36	9.457	1370	1373	1376	rVV	103554	147510	2.57%	0.436%

LSC Area Percent Report

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Operator : JC/MD
Sample : N5614-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

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\82X102622W.M
Title : SW846 8260

37	9.500	1376	1380	1387	rVB4	65717	130383	2.27%	0.385%
38	9.762	1416	1423	1430	rBV3	55100	96167	1.68%	0.284%
39	10.055	1467	1471	1476	rVB	335641	426887	7.44%	1.261%
40	10.195	1489	1494	1502	rVB	980415	1235695	21.52%	3.649%
41	10.305	1508	1512	1517	rVB	155800	215326	3.75%	0.636%
42	10.963	1615	1620	1627	rBV	423221	540000	9.41%	1.595%
43	11.079	1635	1639	1644	rBV	321585	412342	7.18%	1.218%
44	11.305	1671	1676	1681	rBV	345989	407833	7.10%	1.204%
45	11.402	1685	1692	1696	rBV	97481	125087	2.18%	0.369%
46	11.616	1721	1727	1736	rBV	203453	268870	4.68%	0.794%
47	11.866	1763	1768	1770	rBV	61562	83775	1.46%	0.247%
48	11.890	1770	1772	1777	rVB	80877	87825	1.53%	0.259%
49	12.024	1788	1794	1798	rBV	361158	478312	8.33%	1.412%
50	12.244	1824	1830	1836	rBV	4798263	5740871	100.00%	16.952%
51	12.311	1836	1841	1852	rVB2	387104	549298	9.57%	1.622%
52	12.408	1852	1857	1862	rVB2	184776	247160	4.31%	0.730%
53	12.463	1862	1866	1872	rVB3	61048	88720	1.55%	0.262%
54	12.530	1872	1877	1883	rBV	52208	64969	1.13%	0.192%
55	12.615	1883	1891	1893	rBV	282355	353200	6.15%	1.043%
56	12.646	1893	1896	1900	rVB	411959	466816	8.13%	1.378%
57	12.701	1900	1905	1912	rBV	1572074	2073312	36.11%	6.122%
58	12.841	1923	1928	1933	rVB	56083	77538	1.35%	0.229%
59	12.920	1933	1941	1947	rBV	573753	720044	12.54%	2.126%
60	13.024	1953	1958	1966	rVB2	140897	211790	3.69%	0.625%
61	13.115	1966	1973	1974	rBV2	49046	66283	1.15%	0.196%
62	13.134	1974	1976	1978	rVV	95836	114061	1.99%	0.337%
63	13.170	1978	1982	1991	rVB	839354	1056286	18.40%	3.119%
64	13.292	1997	2002	2007	rVB	2283072	2732291	47.59%	8.068%
65	13.341	2007	2010	2017	rVV4	104305	139105	2.42%	0.411%
66	13.426	2020	2024	2031	rVB	259649	351401	6.12%	1.038%
67	13.506	2031	2037	2039	rBV2	107381	144403	2.52%	0.426%
68	13.542	2039	2043	2046	rVV	358422	490198	8.54%	1.447%
69	13.585	2046	2050	2054	rVV2	272333	425338	7.41%	1.256%
70	13.640	2054	2059	2060	rVV2	184164	255629	4.45%	0.755%
71	13.664	2060	2063	2069	rVB2	325562	487786	8.50%	1.440%
72	13.774	2074	2081	2091	rBV	869546	1132604	19.73%	3.344%
73	13.853	2091	2094	2099	rVB	49078	65700	1.14%	0.194%
74	13.926	2099	2106	2110	rBV2	174737	252875	4.40%	0.747%
75	13.969	2110	2113	2117	rVB	148862	168363	2.93%	0.497%
76	14.073	2126	2130	2137	rBV	284648	380035	6.62%	1.122%

LSC Area Percent Report

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Sample : N5614-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

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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Title : SW846 8260

77	14.219	2148	2154	2158	rBV	137820	225884	3.93%	0.667%
78	14.322	2167	2171	2176	rBV	130156	187804	3.27%	0.555%
79	14.371	2176	2179	2184	rVB	195863	235863	4.11%	0.696%
80	14.481	2192	2197	2201	rBV2	65679	96226	1.68%	0.284%
81	14.609	2210	2218	2219	rBV2	42270	75400	1.31%	0.223%
82	14.639	2219	2223	2227	rVB	143137	188566	3.28%	0.557%
83	14.780	2241	2246	2251	rBV	248563	326609	5.69%	0.964%

Sum of corrected areas: 33865290

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111522\
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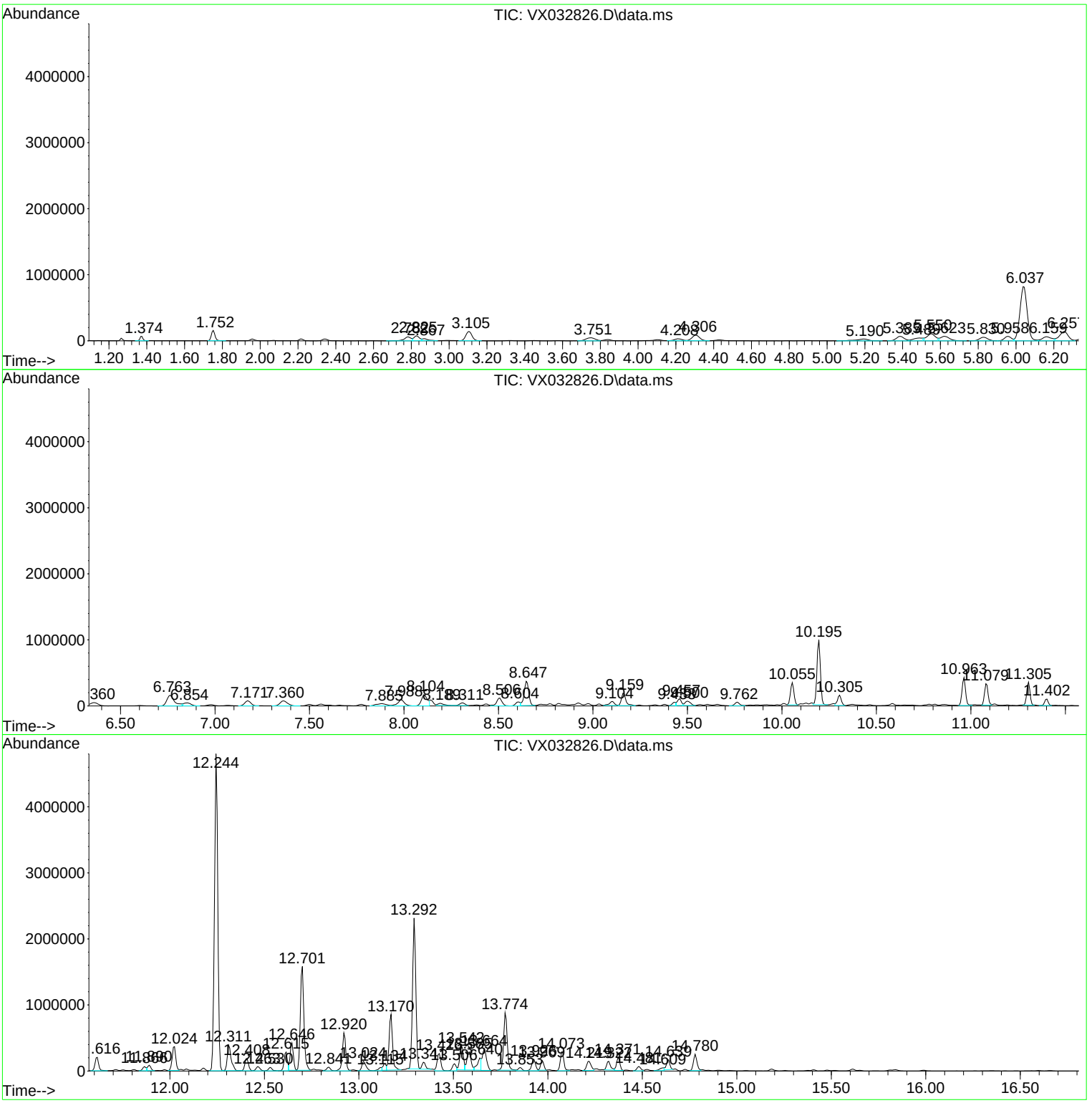
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111522\
Data File : VX032826.D
Acq On : 15 Nov 2022 16:35
Operator : JC/MD
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

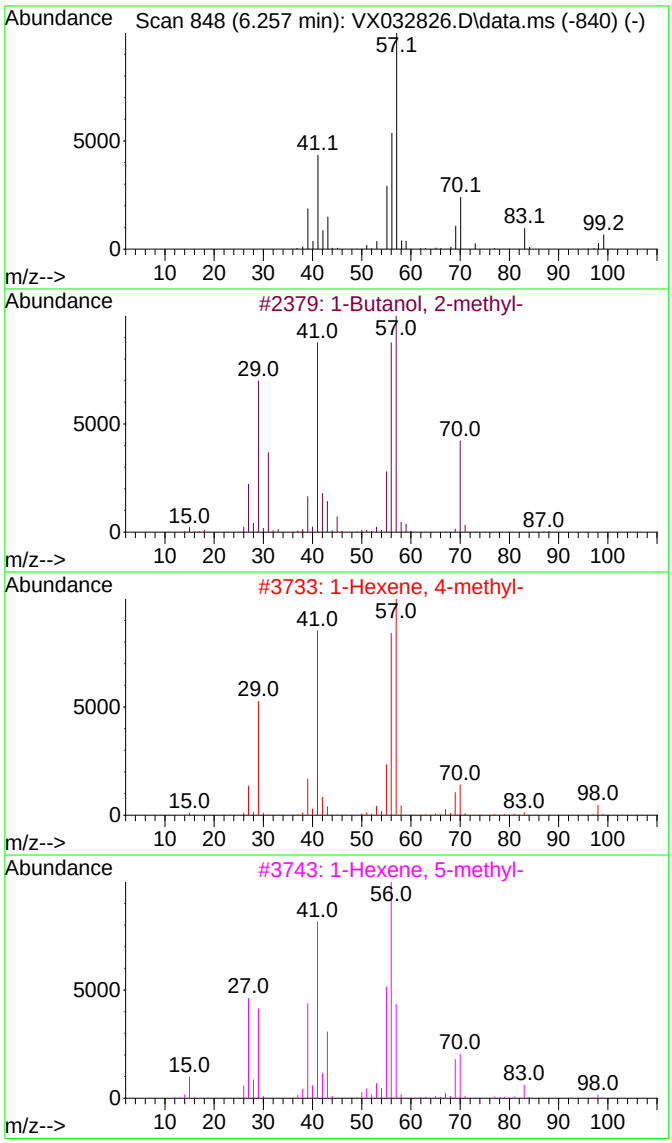
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

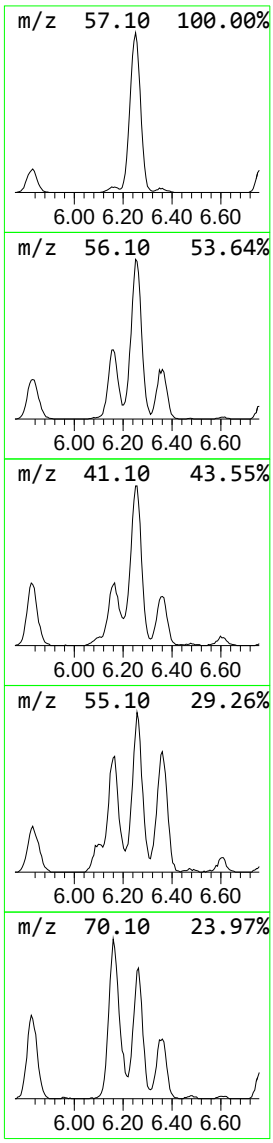
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 1-Butanol, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.257	48.88 ug/l	438138	1,4-Difluorobenzene	6.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Butanol, 2-methyl-	88	C5H12O	000137-32-6	50
2	1-Hexene, 4-methyl-	98	C7H14	003769-23-1	47
3	1-Hexene, 5-methyl-	98	C7H14	003524-73-0	43
4	Acetamide, N-2-propenyl-	99	C5H9NO	000692-33-1	43
5	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	37





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111522\
Data File : VX032826.D
Acq On : 15 Nov 2022 16:35
Operator : JC/MD
Sample : N5614-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

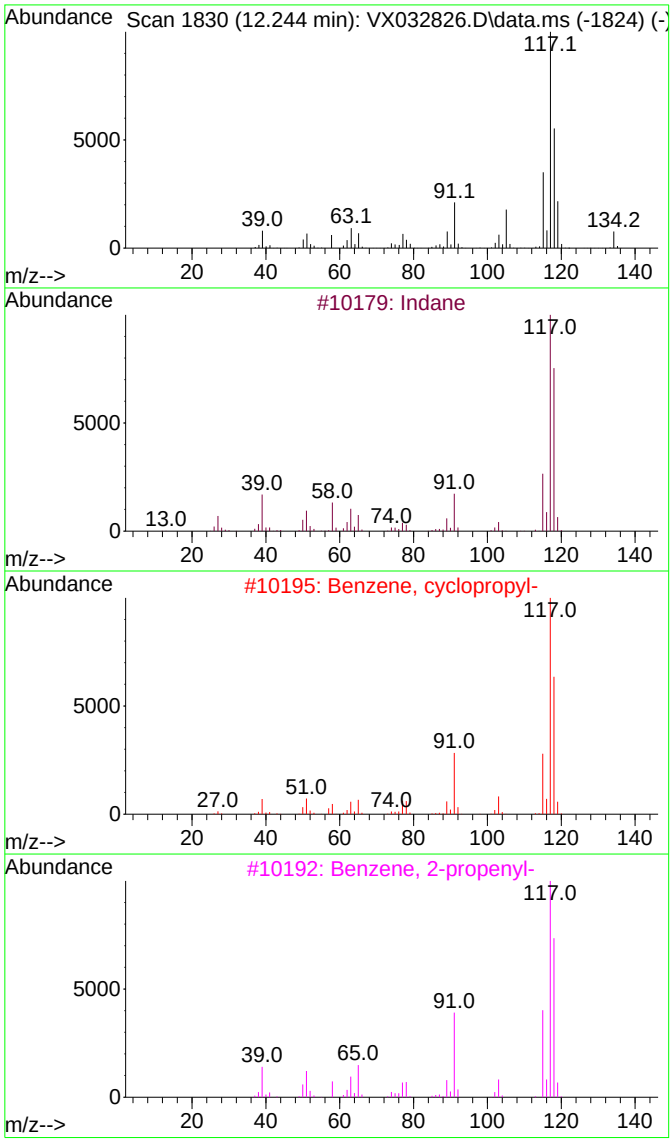
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

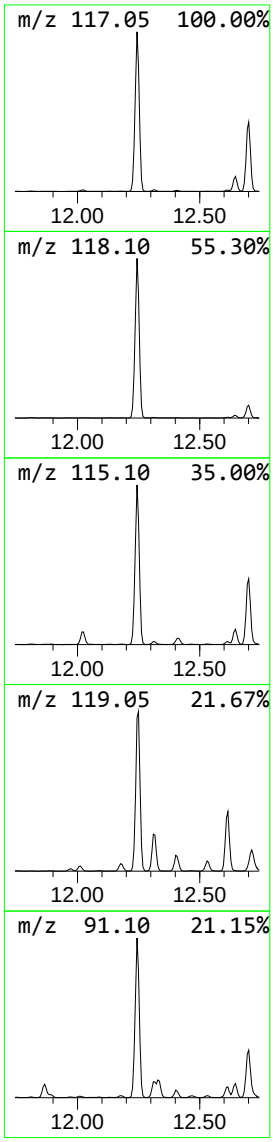
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	600.12 ug/l	5740870	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	76
2	Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	70
4	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	58
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	55





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111522\
Data File : VX032826.D
Acq On : 15 Nov 2022 16:35
Operator : JC/MD
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ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

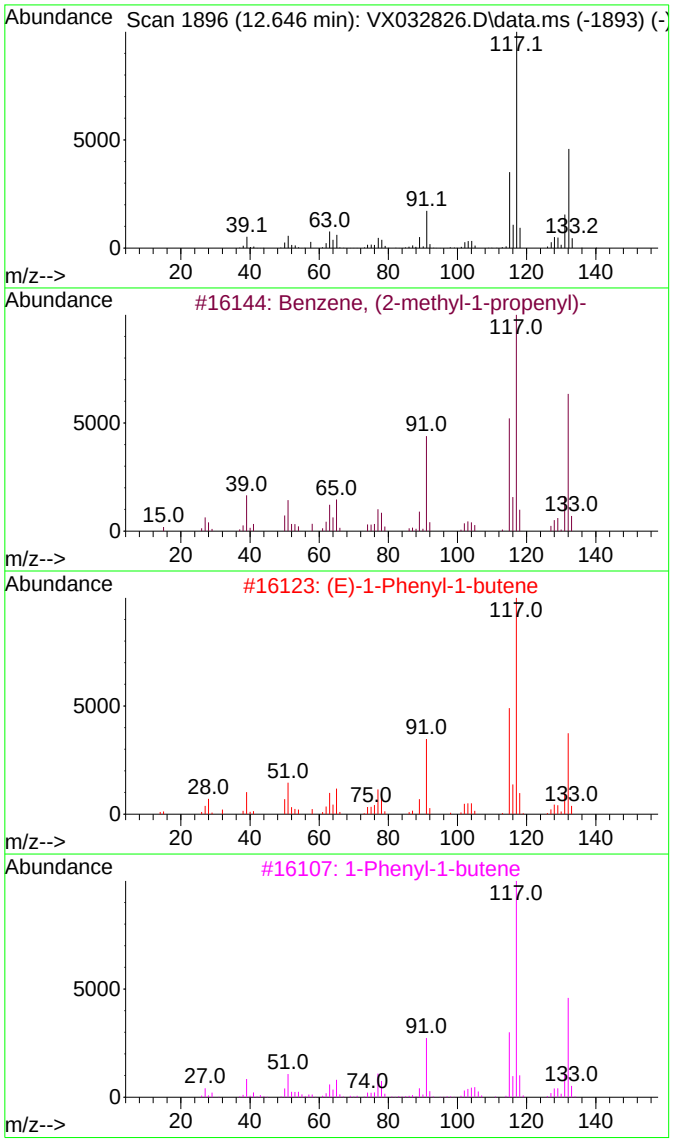
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

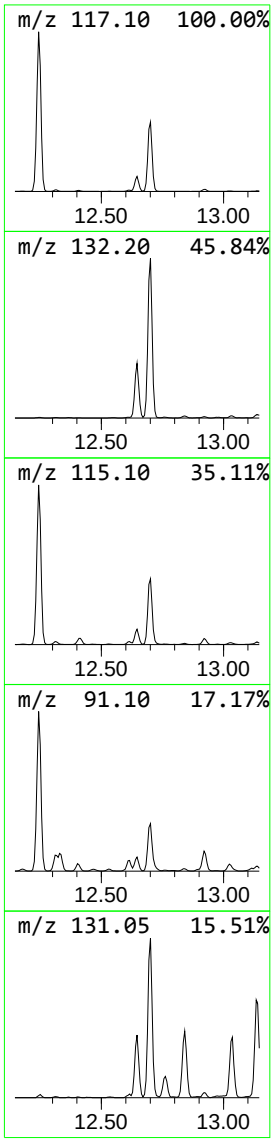
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, (2-methyl-1-propen... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.646	48.80 ug/l	466816	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	94
2	(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	93
3	1-Phenyl-1-butene	132	C10H12	000824-90-8	93
4	Benzene, (1-methylenepropyl)-	132	C10H12	002039-93-2	91
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111522\
Data File : VX032826.D
Acq On : 15 Nov 2022 16:35
Operator : JC/MD
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ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

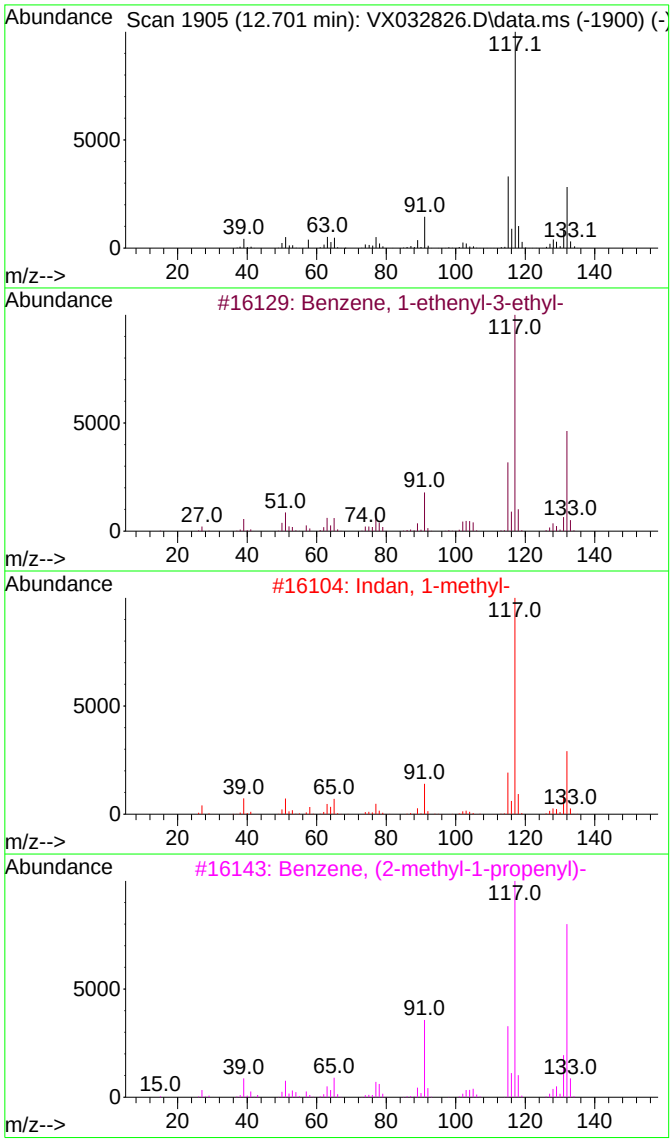
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

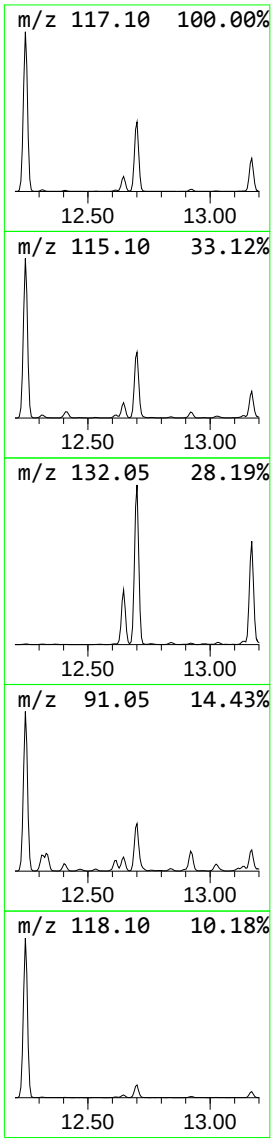
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	216.73 ug/l	2073310	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, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
4	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	83





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111522\
Data File : VX032826.D
Acq On : 15 Nov 2022 16:35
Operator : JC/MD
Sample : N5614-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

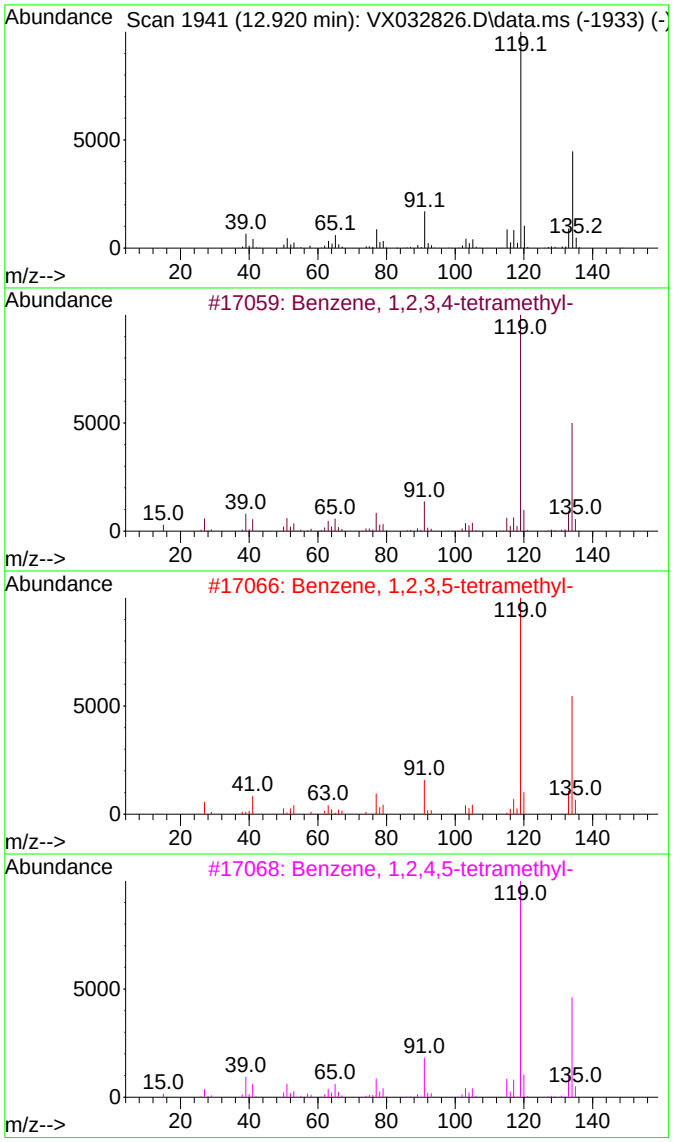
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

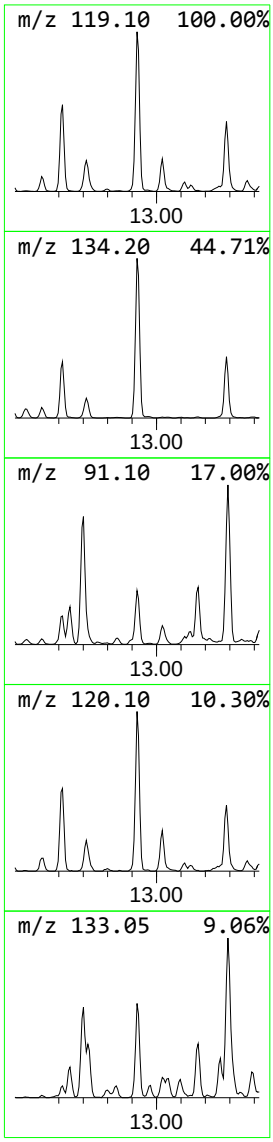
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.920	75.27 ug/l	720044	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	97
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
4	o-Cymene	134	C10H14	000527-84-4	95
5	1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111522\
Data File : VX032826.D
Acq On : 15 Nov 2022 16:35
Operator : JC/MD
Sample : N5614-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

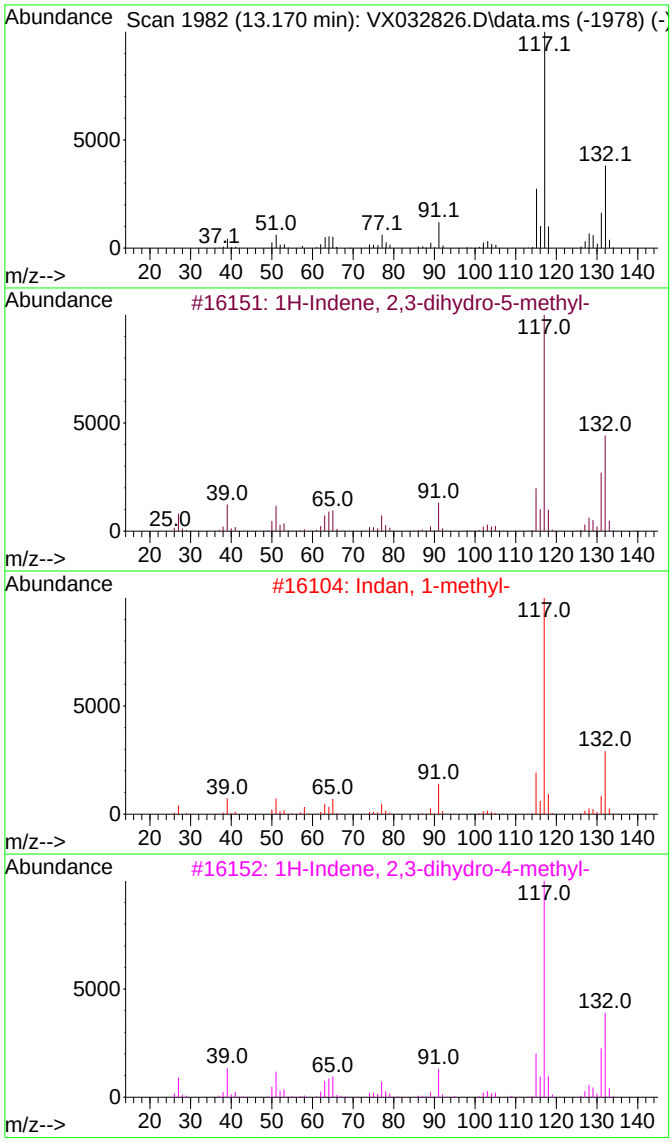
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

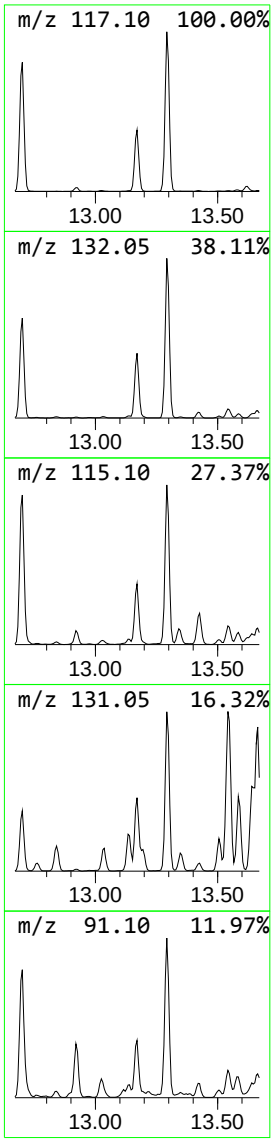
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.170	110.42 ug/l	1056290	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	93
2	Indan, 1-methyl-	132	C10H12	000767-58-8	91
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
4	Benzene, 2-butenyl-	132	C10H12	001560-06-1	87
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111522\
Data File : VX032826.D
Acq On : 15 Nov 2022 16:35
Operator : JC/MD
Sample : N5614-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

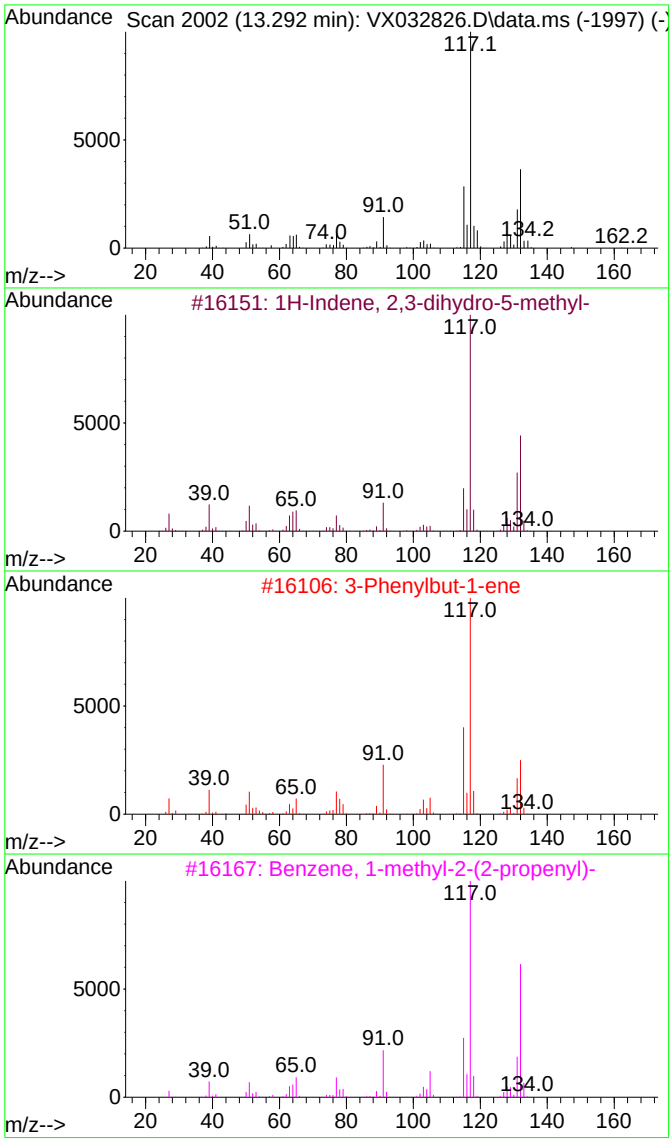
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

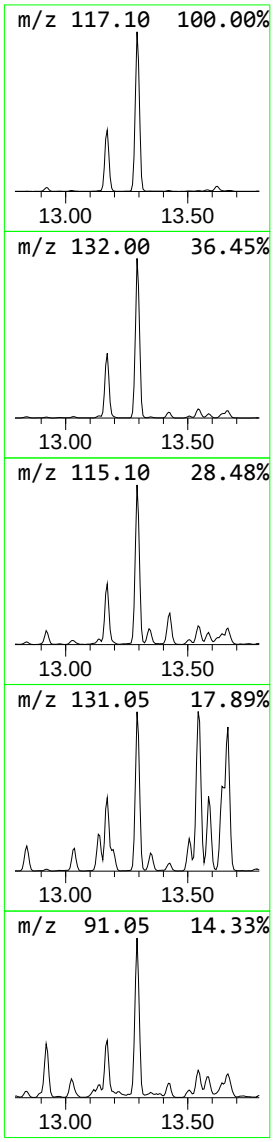
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 3-Phenylbut-1-ene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	285.62 ug/l	2732290	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	91
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
3	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87
4	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
5	1-Phenyl-1-butene	132	C10H12	000824-90-8	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111522\
Data File : VX032826.D
Acq On : 15 Nov 2022 16:35
Operator : JC/MD
Sample : N5614-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

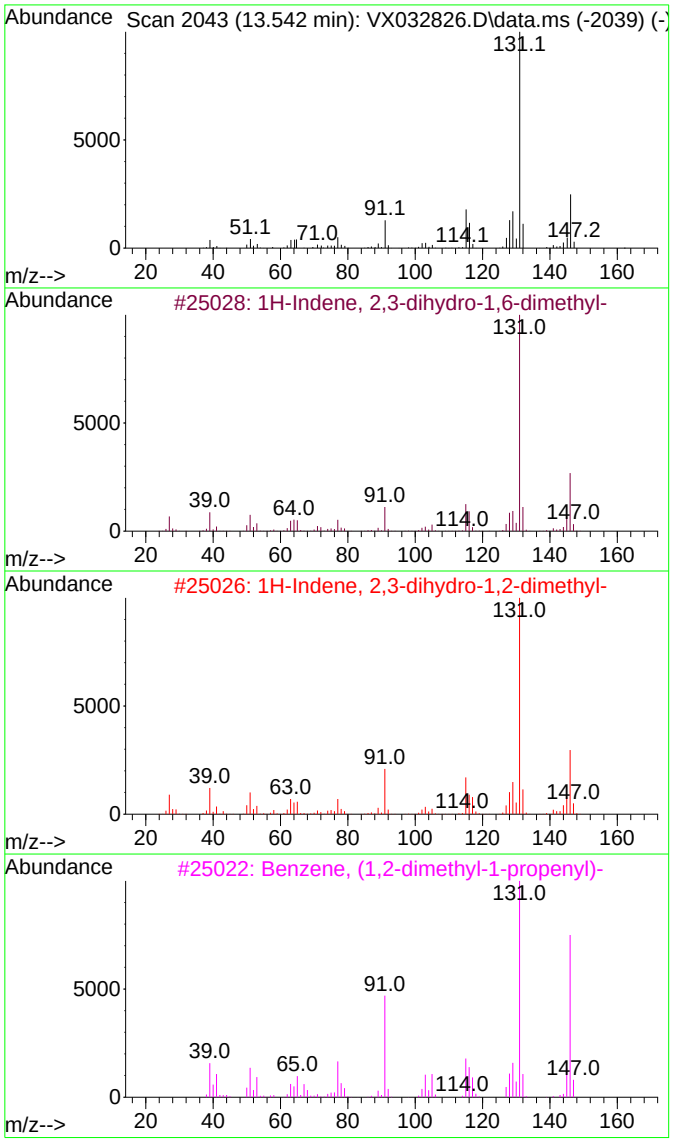
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

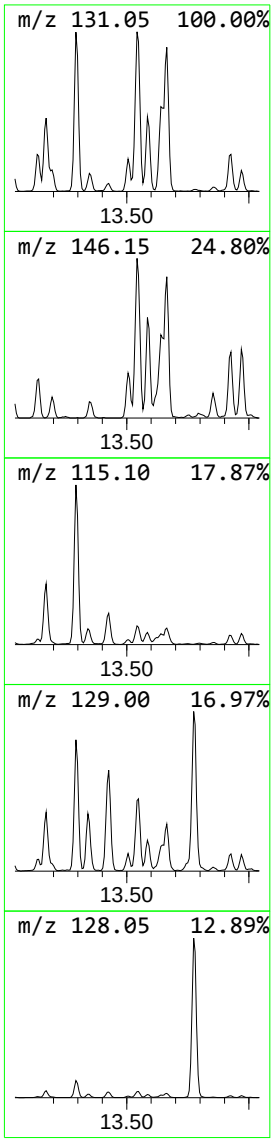
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.542	51.24 ug/l	490198	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	94
2	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	94
3	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	94
4	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	94
5	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111522\
Data File : VX032826.D
Acq On : 15 Nov 2022 16:35
Operator : JC/MD
Sample : N5614-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

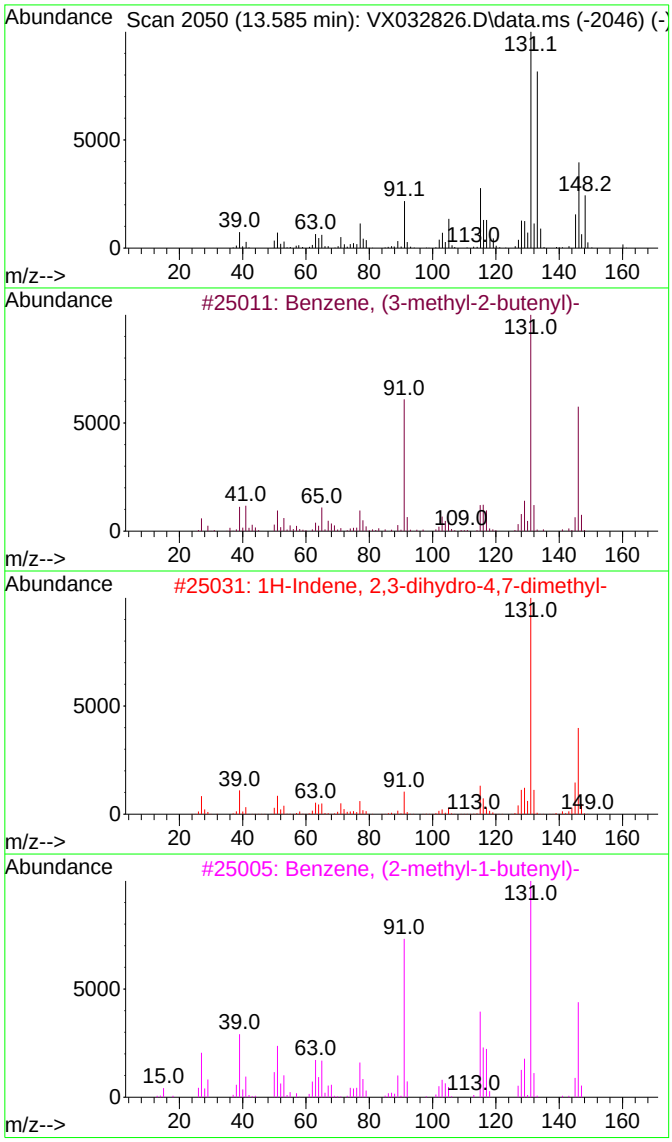
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

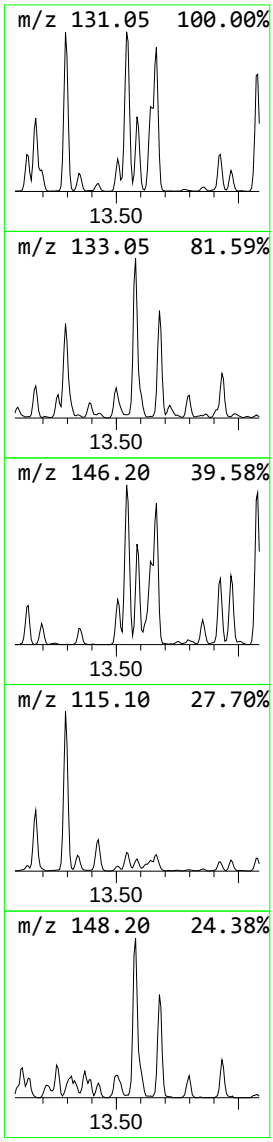
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, (3-methyl-2-butenyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.585	44.46 ug/l	425338	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	91
2	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	90
3	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	83
4	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	83
5	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111522\
Data File : VX032826.D
Acq On : 15 Nov 2022 16:35
Operator : JC/MD
Sample : N5614-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

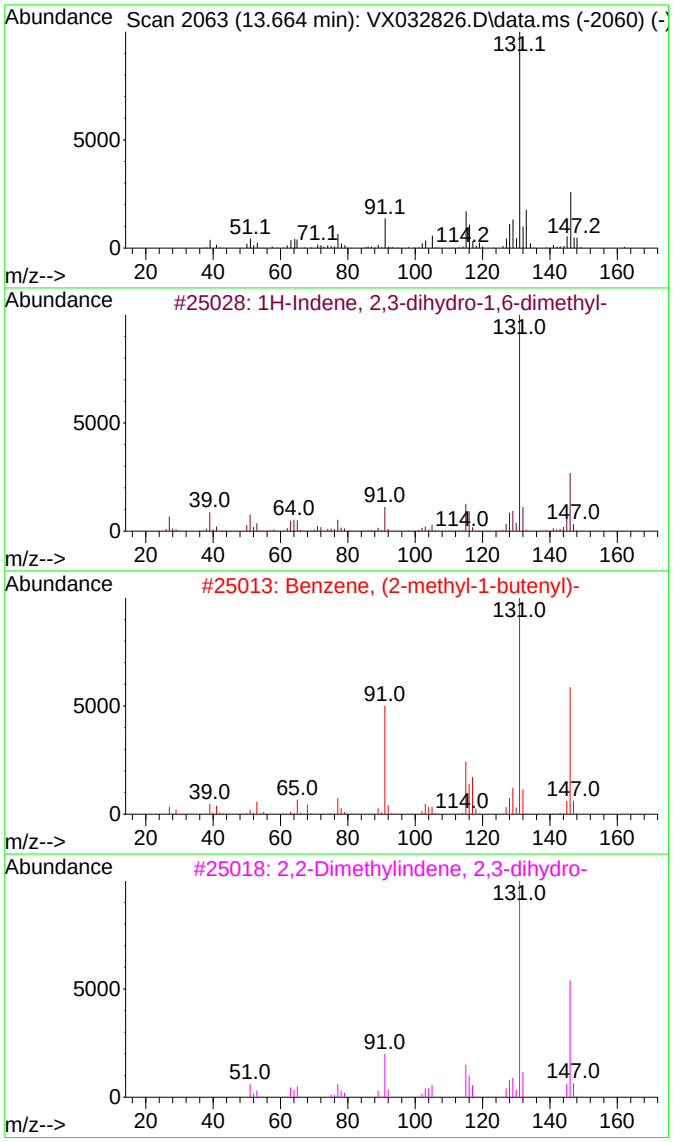
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

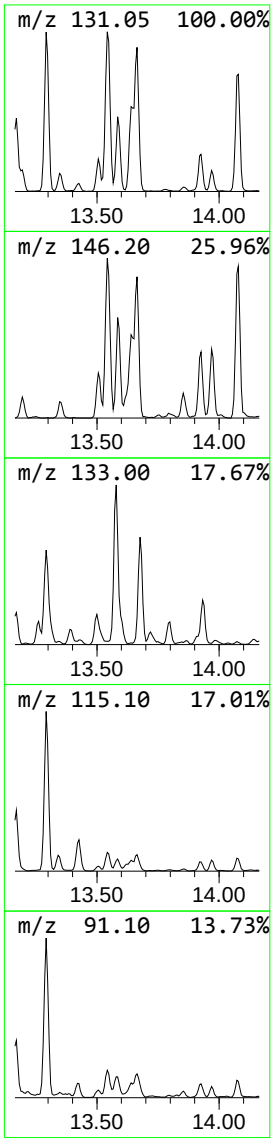
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, (2-methyl-1-butenyl)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.664	50.99 ug/l	487786	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	93
2	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	93
3	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	93
4	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	93
5	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111522\
Data File : VX032826.D
Acq On : 15 Nov 2022 16:35
Operator : JC/MD
Sample : N5614-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.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
1-Butanol, 2-me...	6.257	48.9	ug/l	438138	2	6.763	448213	50.0
Indane	12.244	600.1	ug/l	5740870	4	12.024	478312	50.0
Benzene, (2-met...	12.646	48.8	ug/l	466816	4	12.024	478312	50.0
Benzene, 1-ethe...	12.701	216.7	ug/l	2073310	4	12.024	478312	50.0
Benzene, 1,2,3,...	12.920	75.3	ug/l	720044	4	12.024	478312	50.0
1H-Indene, 2,3-...	13.170	110.4	ug/l	1056290	4	12.024	478312	50.0
3-Phenylbut-1-ene	13.292	285.6	ug/l	2732290	4	12.024	478312	50.0
1H-Indene, 2,3-...	13.542	51.2	ug/l	490198	4	12.024	478312	50.0
Benzene, (3-met...	13.585	44.5	ug/l	425338	4	12.024	478312	50.0
Benzene, (2-met...	13.664	51.0	ug/l	487786	4	12.024	478312	50.0