LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070124\
Data File : VX042116.D
Acq On : 01 Jul 2024 15:55
Operator : JC/MD
Sample : P3092-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW6-20240626

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\82X061824W.M
Title : SW846 8260

Signal : TIC: VX042116.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.240	21	25	27	rBV	58128	56479	1.47%	0.157%
2	1.374	41	47	53	rBV	72440	88125	2.29%	0.245%
3	1.715	98	103	115	rVB	230944	315604	8.19%	0.876%
4	1.935	134	139	148	rVB2	77973	124474	3.23%	0.346%
5	2.197	178	182	192	rVB	42042	64463	1.67%	0.179%
6	2.764	259	275	278	rBV2	47182	138922	3.61%	0.386%
7	2.813	278	283	287	rVV	130273	275486	7.15%	0.765%
8	2.855	287	290	308	rVB2	69276	153792	3.99%	0.427%
9	3.093	319	329	346	rBV3	112655	349205	9.07%	0.970%
10	3.434	377	385	398	rVB	43298	95924	2.49%	0.266%
11	4.294	517	526	540	rVB	107857	309598	8.04%	0.860%
12	5.385	690	705	712	rBV2	78060	240779	6.25%	0.669%
13	5.471	712	719	725	rVV4	35900	121020	3.14%	0.336%
14	5.544	725	731	739	rVB	96330	256113	6.65%	0.711%
15	5.818	765	776	789	rBV2	54566	177385	4.60%	0.493%
16	5.952	790	798	803	rBV	73387	188776	4.90%	0.524%
17	6.032	803	811	824	rVB	1451886	3852177	100.00%	10.698%
18	6.153	824	831	833	rBV2	24493	48924	1.27%	0.136%
19	6.251	841	847	852	rBV2	36924	89872	2.33%	0.250%
20	6.330	852	860	880	rVB	369612	1111802	28.86%	3.087%
21	6.605	892	905	922	rBV6	19628	67094	1.74%	0.186%
22	6.757	922	930	940	rVV	195496	491166	12.75%	1.364%
23	6.836	940	943	955	rVB4	17763	47099	1.22%	0.131%
24	7.373	1016	1031	1044	rBV4	152216	409953	10.64%	1.138%
25	7.653	1071	1077	1090	rVB2	37440	78047	2.03%	0.217%
26	7.982	1122	1131	1143	rVB2	40501	105070	2.73%	0.292%
27	8.104	1143	1151	1160	rBV2	56980	136927	3.55%	0.380%
28	8.446	1200	1207	1212	rBV3	33493	89767	2.33%	0.249%
29	8.525	1212	1220	1226	rVB2	90588	203349	5.28%	0.565%
30	8.598	1226	1232	1235	rBV4	47316	85026	2.21%	0.236%
31	8.647	1235	1240	1246	rVB	422252	662660	17.20%	1.840%
32	8.720	1246	1252	1258	rBV	753414	1188150	30.84%	3.299%
33	8.866	1270	1276	1281	rVB	43068	90397	2.35%	0.251%
34	9.098	1308	1314	1321	rVB	38328	67514	1.75%	0.187%
35	9.458	1365	1373	1376	rBV3	21755	39898	1.04%	0.111%
36	9.507	1376	1381	1385	rVV5	27152	57163	1.48%	0.159%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070124\
Data File : VX042116.D
Acq On : 01 Jul 2024 15:55
Operator : JC/MD
Sample : P3092-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW6-20240626

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\82X061824W.M
Title : SW846 8260

37	9.555	1385	1389	1395	rVB3	39581	71038	1.84%	0.197%
38	10.055	1460	1471	1475	rBV	457905	645341	16.75%	1.792%
39	10.195	1488	1494	1504	rBV	1513369	2015017	52.31%	5.596%
40	10.299	1504	1511	1518	rVB	2836287	3796209	98.55%	10.542%
41	10.640	1562	1567	1580	rVB	2068900	2605070	67.63%	7.234%
42	10.964	1613	1620	1626	rVB	116474	169171	4.39%	0.470%
43	11.079	1634	1639	1645	rBV	416834	532945	13.83%	1.480%
44	11.305	1672	1676	1683	rVB	146947	181962	4.72%	0.505%
45	11.378	1683	1688	1690	rBV	537488	730099	18.95%	2.027%
46	11.451	1695	1700	1711	rVB	643967	805020	20.90%	2.236%
47	11.610	1721	1726	1737	rBV	1024026	1257318	32.64%	3.492%
48	11.750	1744	1749	1756	rVB	2628995	3165561	82.18%	8.791%
49	11.969	1781	1785	1788	rBV	54307	63339	1.64%	0.176%
50	12.018	1788	1793	1799	rVB	522328	678368	17.61%	1.884%
51	12.085	1799	1804	1812	rBV	969080	1147381	29.79%	3.186%
52	12.244	1823	1830	1838	rBV3	797225	1244602	32.31%	3.456%
53	12.317	1838	1842	1851	rVB2	328992	466343	12.11%	1.295%
54	12.408	1852	1857	1861	rBV2	45754	63279	1.64%	0.176%
55	12.463	1861	1866	1872	rVB	131290	168386	4.37%	0.468%
56	12.530	1872	1877	1879	rBV	232729	298676	7.75%	0.829%
57	12.555	1879	1881	1886	rVB	186392	194882	5.06%	0.541%
58	12.610	1886	1890	1894	rBV	375551	470573	12.22%	1.307%
59	12.646	1894	1896	1900	rVB	69674	69862	1.81%	0.194%
60	12.695	1900	1904	1912	rVB2	263019	395853	10.28%	1.099%
61	12.847	1925	1929	1934	rVB	117555	145455	3.78%	0.404%
62	12.921	1934	1941	1944	rBV	246464	296145	7.69%	0.822%
63	12.963	1944	1948	1954	rVB	303705	376664	9.78%	1.046%
64	13.024	1954	1958	1960	rBV2	31558	47114	1.22%	0.131%
65	13.164	1978	1981	1986	rVB	141123	171976	4.46%	0.478%
66	13.256	1991	1996	1997	rBV2	29470	38577	1.00%	0.107%
67	13.292	1997	2002	2007	rVB2	399405	524442	13.61%	1.456%
68	13.420	2020	2023	2029	rVB	37226	48039	1.25%	0.133%
69	13.500	2031	2036	2039	rBV2	28751	41799	1.09%	0.116%
70	13.542	2039	2043	2046	rVV	53123	68749	1.78%	0.191%
71	13.579	2046	2049	2056	rVV4	74136	145482	3.78%	0.404%
72	13.664	2060	2063	2068	rVB2	64159	95659	2.48%	0.266%
73	13.774	2075	2081	2086	rBV	457690	564451	14.65%	1.567%
74	14.213	2148	2153	2159	rBV	31650	47789	1.24%	0.133%
75	14.634	2218	2222	2227	rVB	120753	148827	3.86%	0.413%
76	14.780	2241	2246	2260	rVB	98929	134390	3.49%	0.373%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070124\
Data File : VX042116.D
Acq On : 01 Jul 2024 15:55
Operator : JC/MD
Sample : P3092-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW6-20240626

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\82X061824W.M
Title : SW846 8260

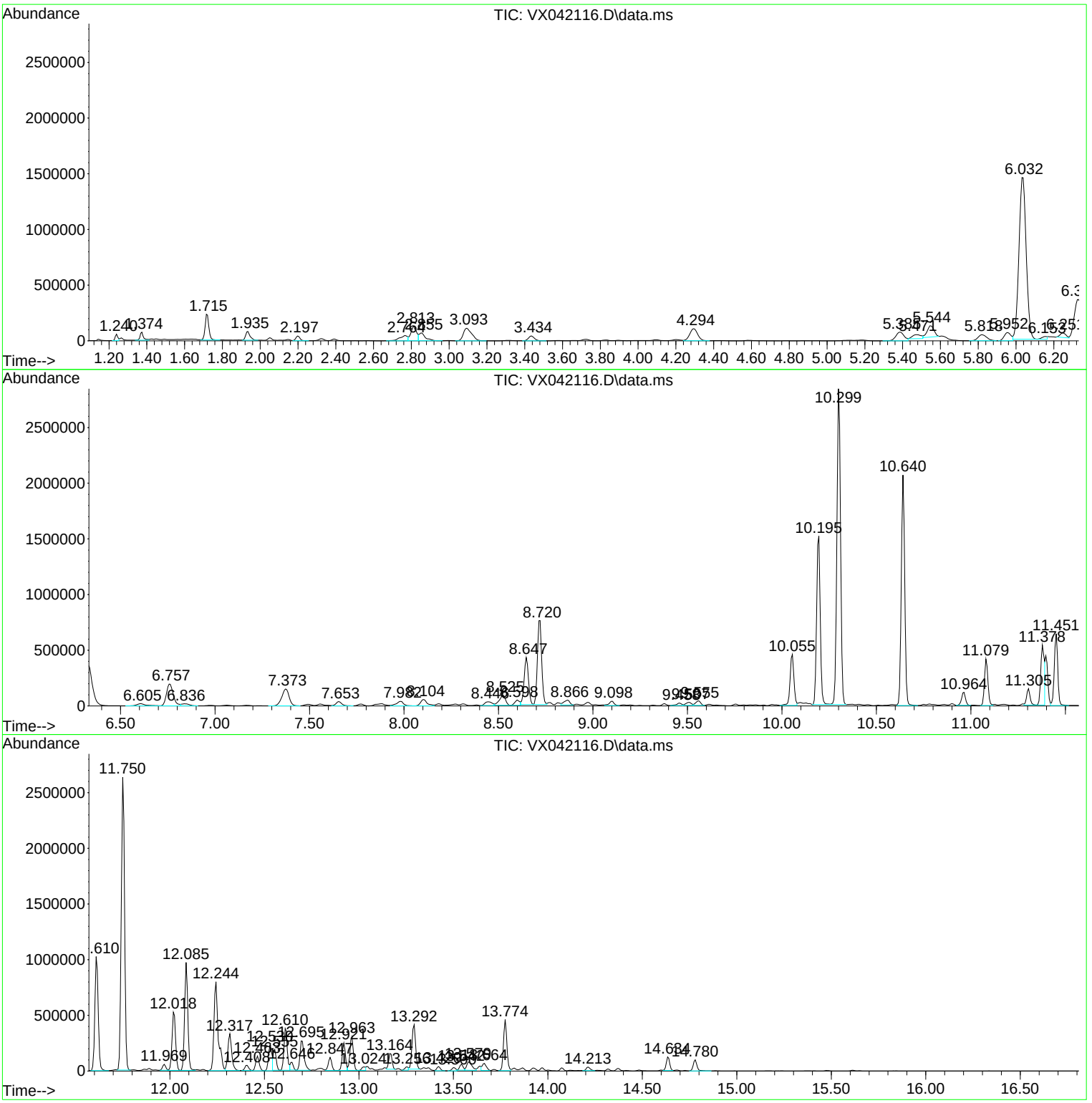
Sum of corrected areas: 36010053

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070124\
Data File : VX042116.D
Acq On : 01 Jul 2024 15:55
Operator : JC/MD
Sample : P3092-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW6-20240626

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061824W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070124\
Data File : VX042116.D
Acq On : 01 Jul 2024 15:55
Operator : JC/MD
Sample : P3092-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW6-20240626

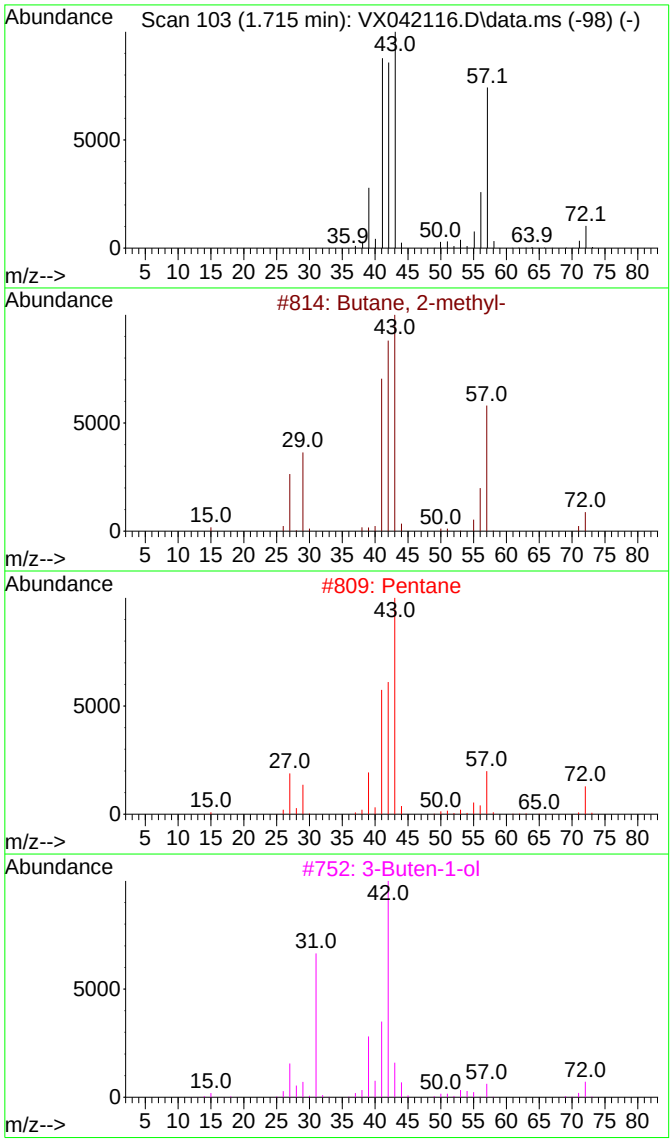
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061824W.M
Quant Title : SW846 8260

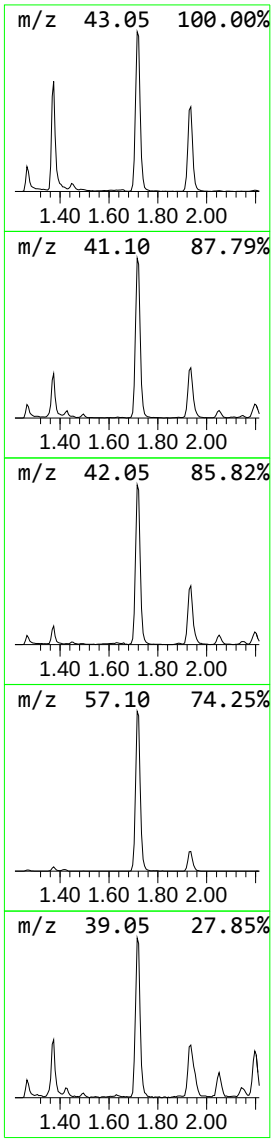
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Butane, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.715	61.61 ug/l	315604	Pentafluorobenzene	5.550

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	90
2	Pentane	72	C5H12	000109-66-0	43
3	3-Buten-1-ol	72	C4H8O	000627-27-0	37
4	Butane, 1-chloro-2-methyl-	106	C5H11Cl	000616-13-7	10
5	1-Butene	56	C4H8	000106-98-9	10





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070124\
Data File : VX042116.D
Acq On : 01 Jul 2024 15:55
Operator : JC/MD
Sample : P3092-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW6-20240626

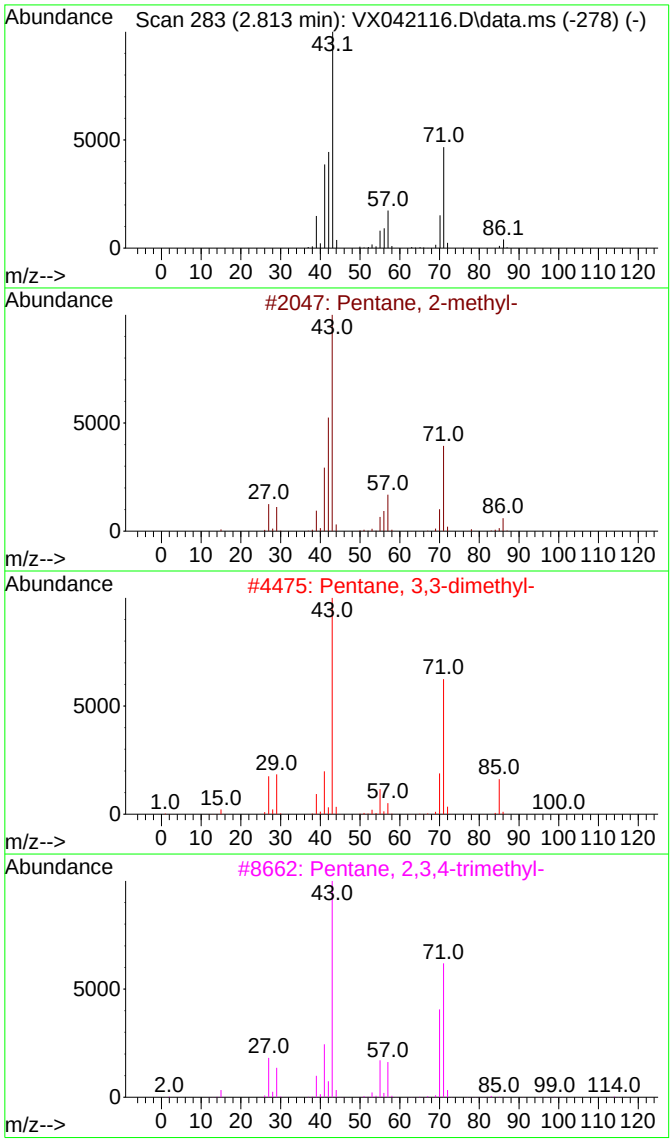
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061824W.M
Quant Title : SW846 8260

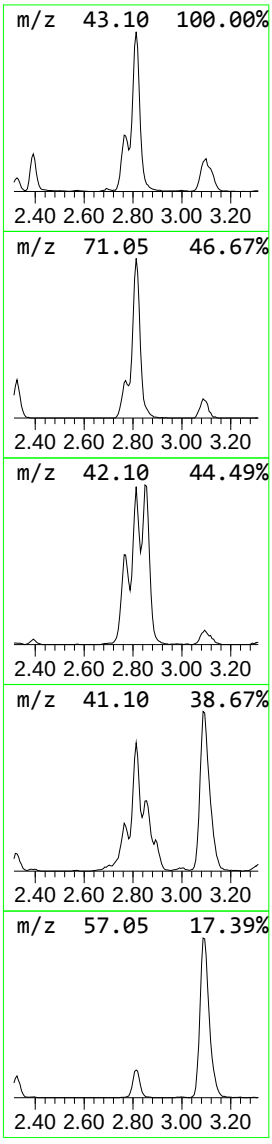
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Pentane, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.813	53.78 ug/l	275486	Pentafluorobenzene	5.550

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2-methyl-	86	C6H14	000107-83-5	91
2	Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	37
3	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	25
4	Heptane, 4-methyl-	114	C8H18	000589-53-7	25
5	Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	25





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070124\
Data File : VX042116.D
Acq On : 01 Jul 2024 15:55
Operator : JC/MD
Sample : P3092-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW6-20240626

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061824W.M
Quant Title : SW846 8260

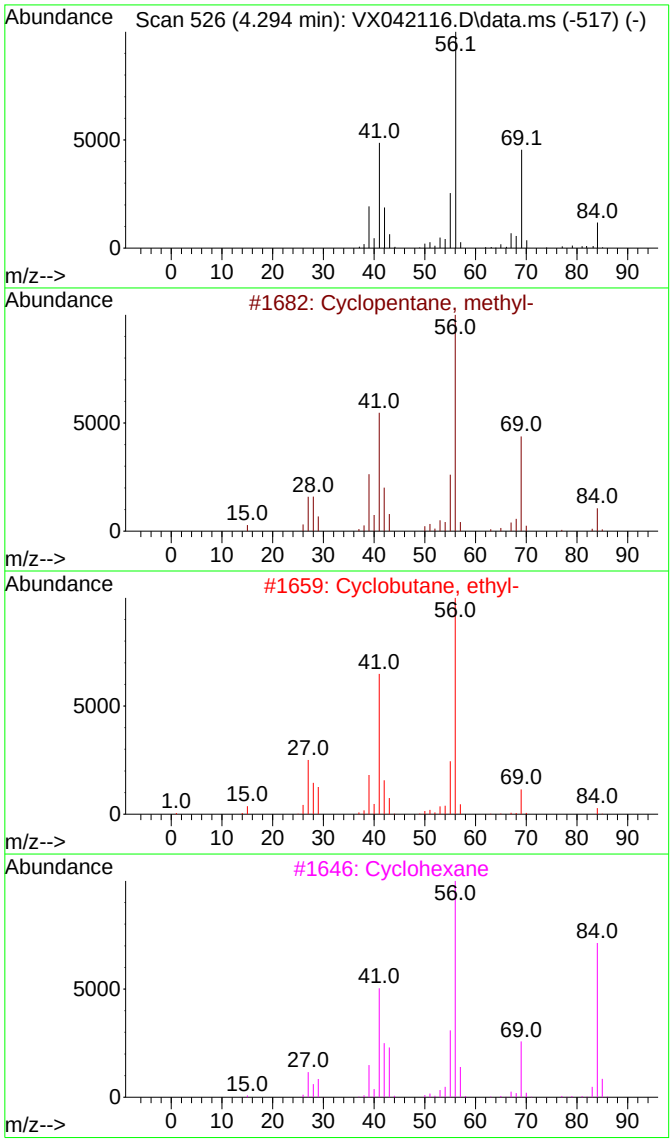
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

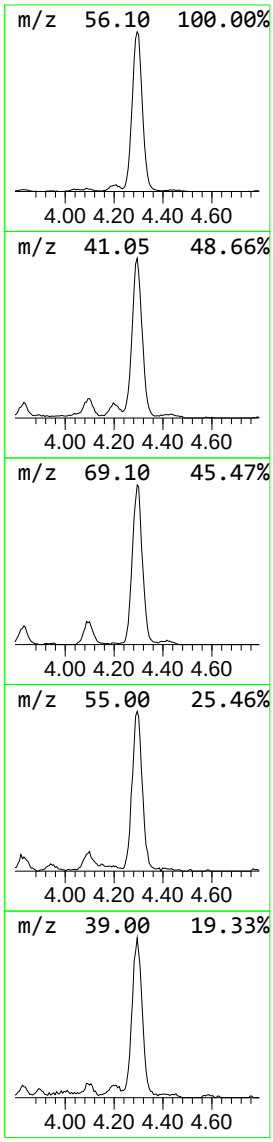
****

Peak Number 3 Cyclopentane, methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.294	60.44 ug/l	309598	Pentafluorobenzene	5.550

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2	Cyclobutane, ethyl-	84	C6H12	004806-61-5	78
3	Cyclohexane	84	C6H12	000110-82-7	74
4	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64
5	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	43





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070124\
Data File : VX042116.D
Acq On : 01 Jul 2024 15:55
Operator : JC/MD
Sample : P3092-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW6-20240626

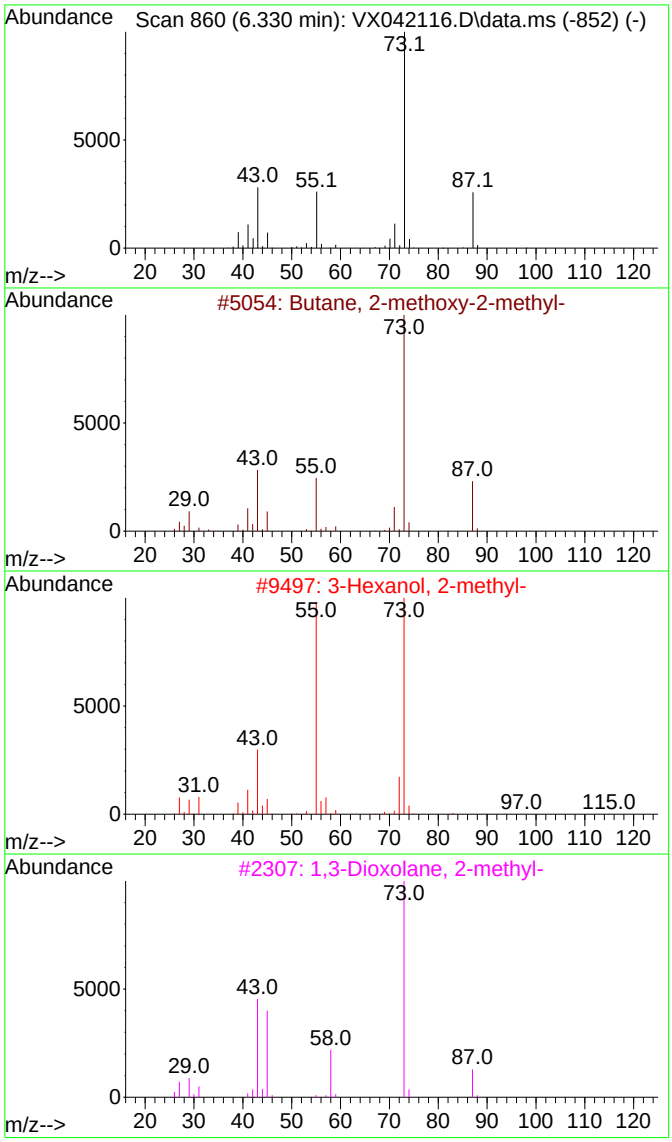
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061824W.M
Quant Title : SW846 8260

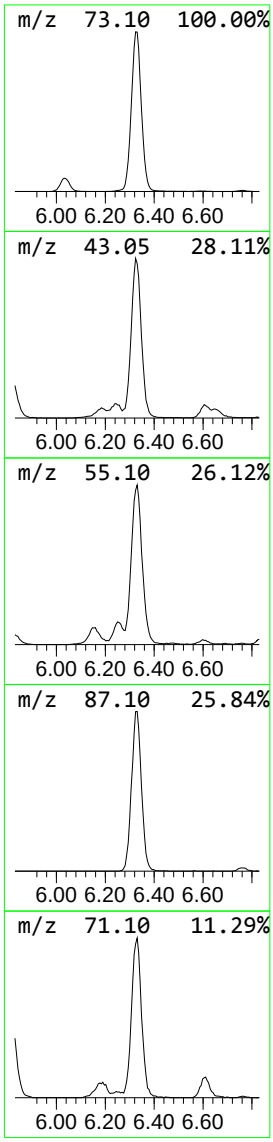
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.330	113.18 ug/l	1111800	1,4-Difluorobenzene	6.757

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
2	3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	17
3	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
4	1,4-Butanediamine	88	C4H12N2	000110-60-1	9
5	Borane, dimethoxy-	74	C2H7BO2	004542-61-4	9





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070124\
Data File : VX042116.D
Acq On : 01 Jul 2024 15:55
Operator : JC/MD
Sample : P3092-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW6-20240626

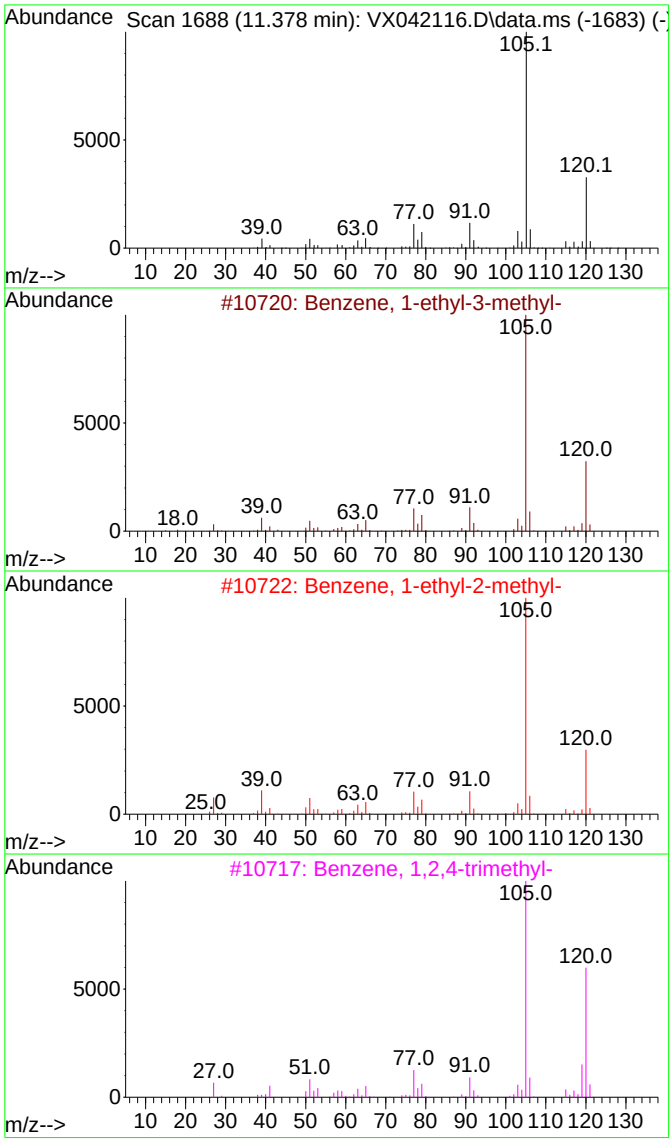
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061824W.M
Quant Title : SW846 8260

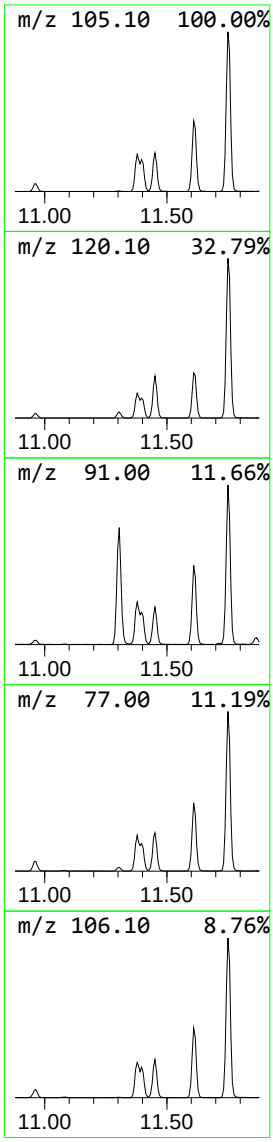
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1-ethyl-4-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.378	53.81 ug/l	730099	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	97
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5	Mesitylene	120	C9H12	000108-67-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070124\
Data File : VX042116.D
Acq On : 01 Jul 2024 15:55
Operator : JC/MD
Sample : P3092-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW6-20240626

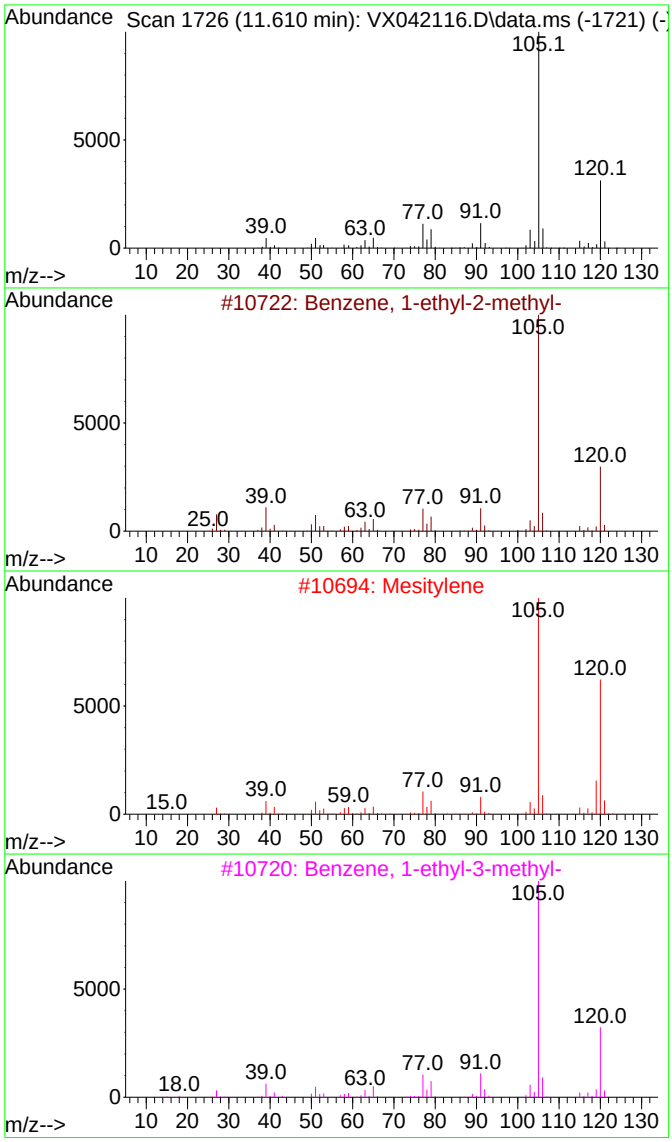
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061824W.M
Quant Title : SW846 8260

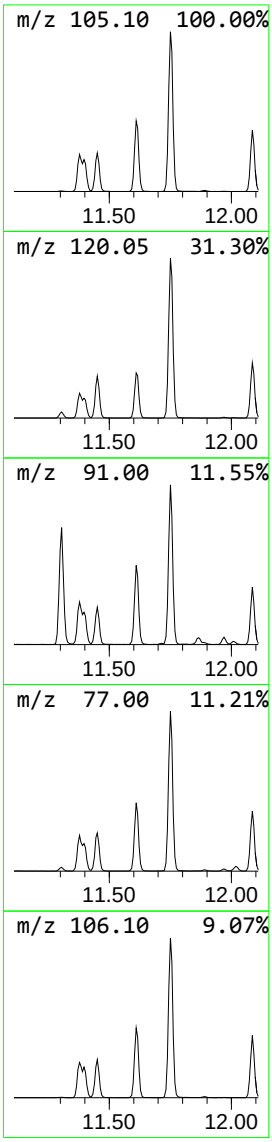
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.610	92.67 ug/l	1257320	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	95
2	Mesitylene	120	C9H12	000108-67-8	91
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070124\
Data File : VX042116.D
Acq On : 01 Jul 2024 15:55
Operator : JC/MD
Sample : P3092-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW6-20240626

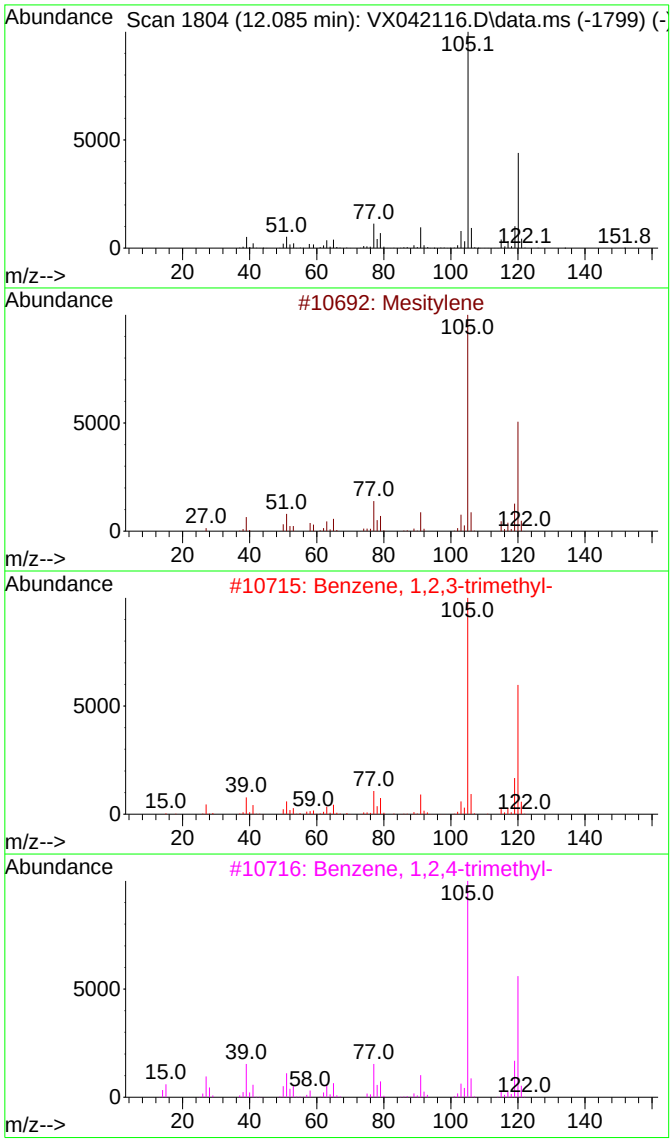
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061824W.M
Quant Title : SW846 8260

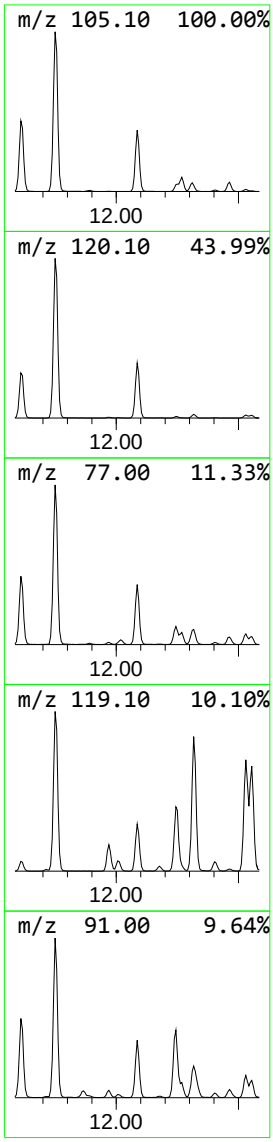
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1-ethyl-3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.085	84.57 ug/l	1147380	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, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070124\
Data File : VX042116.D
Acq On : 01 Jul 2024 15:55
Operator : JC/MD
Sample : P3092-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW6-20240626

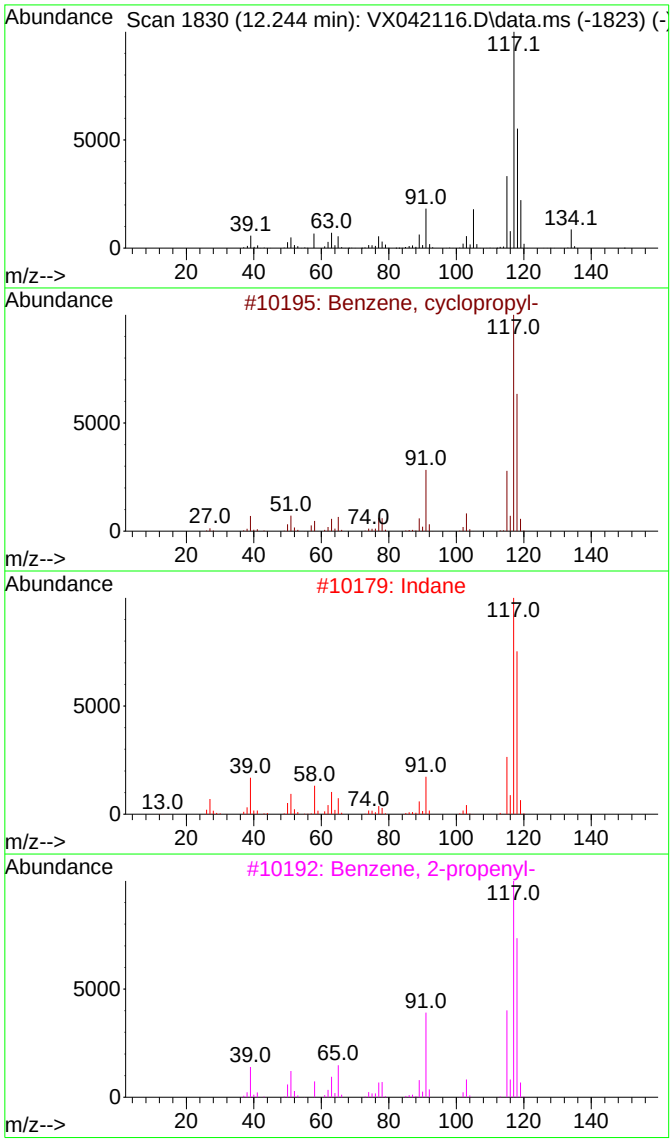
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061824W.M
Quant Title : SW846 8260

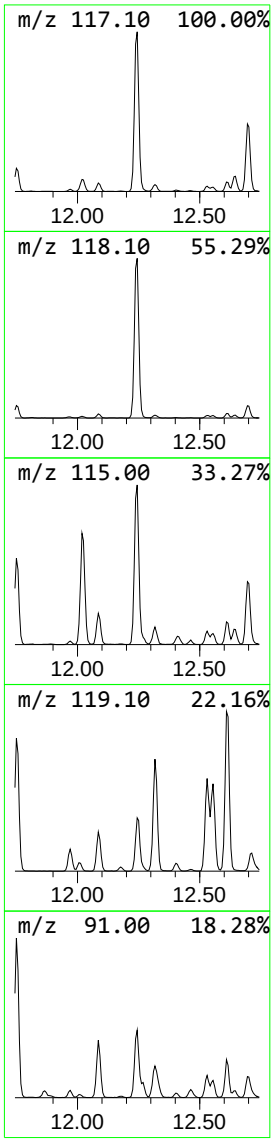
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, cyclopropyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	91.73 ug/l	1244600	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
2	Indane	118	C9H10	000496-11-7	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	58





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070124\
Data File : VX042116.D
Acq On : 01 Jul 2024 15:55
Operator : JC/MD
Sample : P3092-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW6-20240626

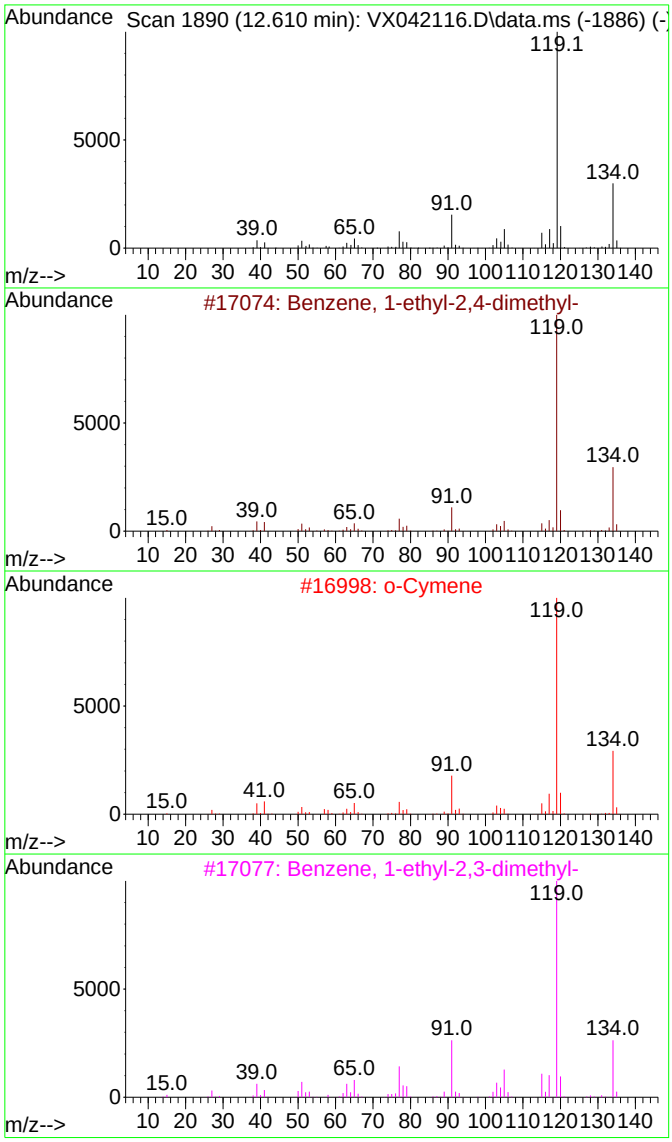
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061824W.M
Quant Title : SW846 8260

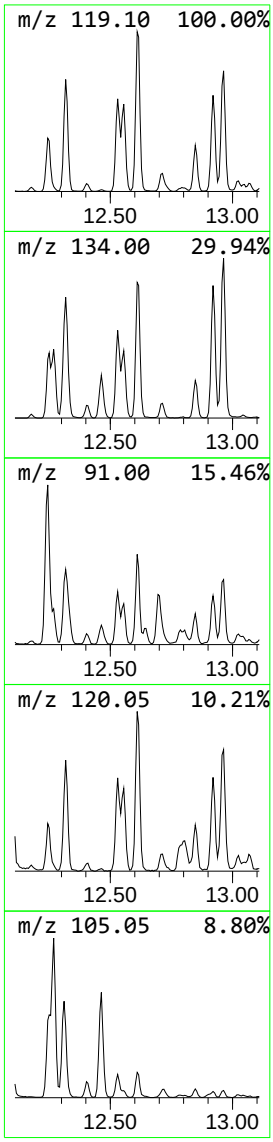
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.610	34.68 ug/l	470573	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2	o-Cymene	134	C10H14	000527-84-4	97
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070124\
Data File : VX042116.D
Acq On : 01 Jul 2024 15:55
Operator : JC/MD
Sample : P3092-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW6-20240626

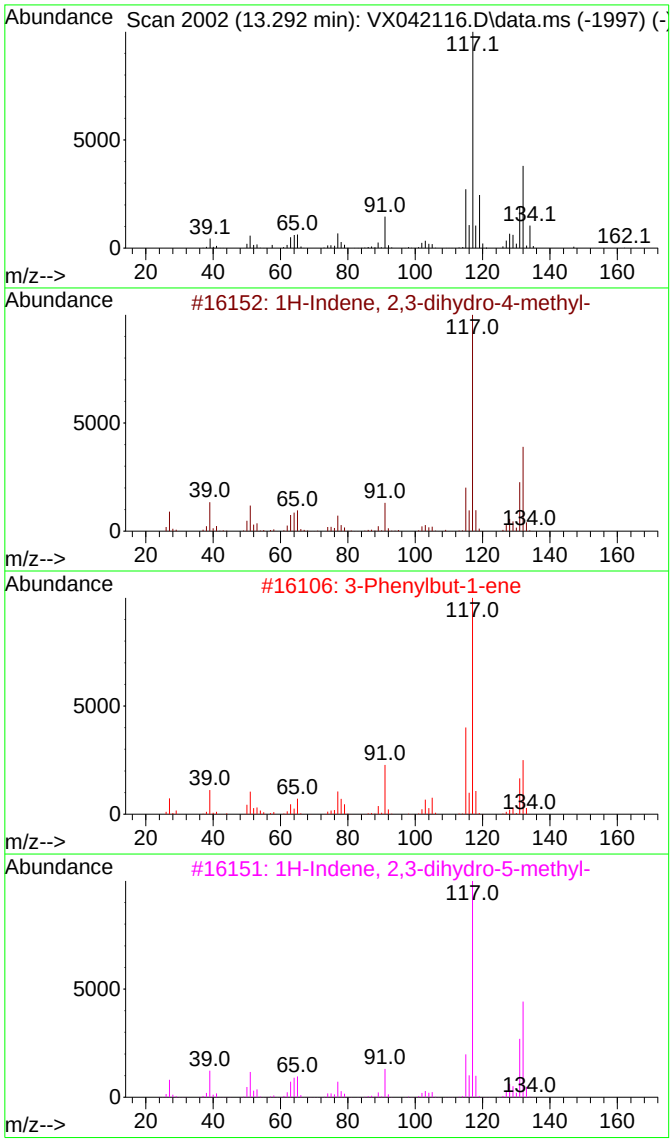
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061824W.M
Quant Title : SW846 8260

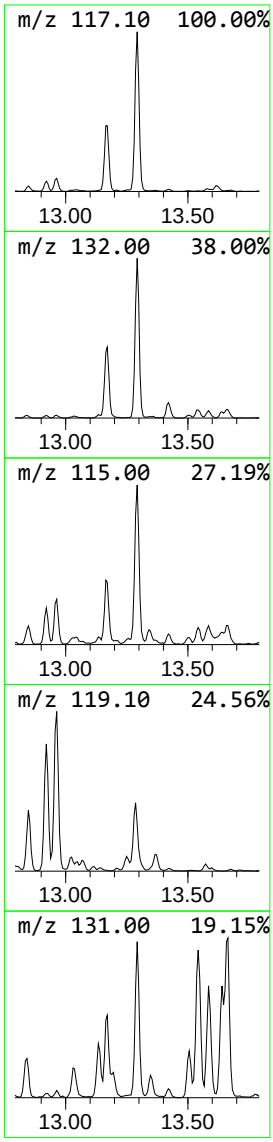
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	38.65 ug/l	524442	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	81
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
3	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	76
4	Benzene, 2-butenyl-	132	C10H12	001560-06-1	74
5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	72





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070124\
Data File : VX042116.D
Acq On : 01 Jul 2024 15:55
Operator : JC/MD
Sample : P3092-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW6-20240626

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061824W.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-	1.715	61.6	ug/l	315604	1	5.550	256113	50.0
Pentane, 2-methyl-	2.813	53.8	ug/l	275486	1	5.550	256113	50.0
Cyclopentane, m...	4.294	60.4	ug/l	309598	1	5.550	256113	50.0
Butane, 2-metho...	6.330	113.2	ug/l	1111800	2	6.757	491166	50.0
Benzene, 1-ethy...	11.378	53.8	ug/l	730099	4	12.024	678368	50.0
Benzene, 1-ethy...	11.610	92.7	ug/l	1257320	4	12.024	678368	50.0
Benzene, 1-ethy...	12.085	84.6	ug/l	1147380	4	12.024	678368	50.0
Benzene, cyclop...	12.244	91.7	ug/l	1244600	4	12.024	678368	50.0
Benzene, 1-ethy...	12.610	34.7	ug/l	470573	4	12.024	678368	50.0
1H-Indene, 2,3-...	13.292	38.7	ug/l	524442	4	12.024	678368	50.0