LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120621\
Data File : VX025563.D
Acq On : 06 Dec 2021 20:49
Operator : JC/MD
Sample : M4940-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6

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\82X111721W.M
Title : SW846 8260

Signal : TIC: VX025563.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.270	22	30	42	rBV2	19501	59497	6.96%	0.754%
2	1.368	43	46	52	rVB	65760	66670	7.80%	0.845%
3	1.746	103	108	117	rVB	206283	285186	33.37%	3.615%
4	1.953	137	142	155	rVB2	80679	121946	14.27%	1.546%
5	2.215	180	185	196	rVB	46208	73888	8.64%	0.937%
6	2.337	199	205	210	rVV2	5027	9646	1.13%	0.122%
7	2.758	260	274	275	rBV2	11628	24647	2.88%	0.312%
8	2.782	275	278	281	rVV	14752	27463	3.21%	0.348%
9	2.825	281	285	288	rVV	32270	63088	7.38%	0.800%
10	2.867	288	292	303	rVV2	42023	95964	11.23%	1.217%
11	2.971	303	309	322	rVB	5648	13055	1.53%	0.165%
12	3.117	322	333	345	rBV4	58024	154185	18.04%	1.955%
13	3.446	380	387	395	rBV2	8843	20548	2.40%	0.260%
14	3.733	426	434	444	rBV2	12787	30771	3.60%	0.390%
15	3.837	444	451	459	rBV2	8069	19685	2.30%	0.250%
16	4.105	486	495	504	rVB3	7806	20108	2.35%	0.255%
17	4.306	519	528	539	rVV2	43086	122619	14.35%	1.554%
18	4.434	539	549	561	rVB3	4058	12584	1.47%	0.160%
19	5.196	665	674	687	rVB2	8164	25348	2.97%	0.321%
20	5.391	694	706	714	rBV2	55975	164409	19.24%	2.084%
21	5.477	714	720	724	rVV3	13727	35853	4.19%	0.454%
22	5.556	725	733	742	rVB	88944	248167	29.03%	3.146%
23	5.830	765	778	788	rBV2	9939	29777	3.48%	0.377%
24	5.964	788	800	805	rVV	59065	154491	18.07%	1.958%
25	6.044	805	813	825	rVV	321105	854734	100.00%	10.835%
26	6.251	826	847	857	rVV5	20215	105898	12.39%	1.342%
27	6.367	858	866	875	rVB7	5112	15283	1.79%	0.194%
28	6.769	922	932	941	rBV	145563	352088	41.19%	4.463%
29	6.848	941	945	955	rVB3	6201	17433	2.04%	0.221%
30	7.171	993	998	1006	rVB4	5872	12072	1.41%	0.153%
31	7.379	1023	1032	1042	rVB4	16576	43987	5.15%	0.558%
32	7.988	1123	1132	1141	rVB	12344	26632	3.12%	0.338%
33	8.110	1145	1152	1155	rBV	18098	35812	4.19%	0.454%
34	8.147	1155	1158	1163	rVV	14797	25818	3.02%	0.327%
35	8.433	1199	1205	1210	rBV3	7788	13698	1.60%	0.174%

LSC Area Percent Report

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Operator : JC/MD
Sample : M4940-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6

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\82X111721W.M
Title : SW846 8260

36	8.507	1212	1217	1225	rVB3	14084	23815	2.79%	0.302%
37	8.653	1225	1241	1247	rBV	309500	488509	57.15%	6.193%
38	8.720	1247	1252	1259	rVB	171400	256977	30.07%	3.258%
39	8.842	1265	1272	1279	rVB3	8094	17219	2.01%	0.218%
40	9.165	1321	1325	1329	rBV	5903	8986	1.05%	0.114%
41	9.458	1365	1373	1377	rBV2	7960	13710	1.60%	0.174%
42	10.055	1466	1471	1477	rVV	297602	393280	46.01%	4.985%
43	10.195	1489	1494	1500	rBV	205230	261135	30.55%	3.310%
44	10.305	1506	1512	1518	rVB	163681	226228	26.47%	2.868%
45	10.646	1563	1568	1576	rVB	61365	79792	9.34%	1.011%
46	10.963	1613	1620	1626	rVB	32572	43227	5.06%	0.548%
47	11.085	1634	1640	1649	rVB	231571	300488	35.16%	3.809%
48	11.305	1667	1676	1684	rBV	55076	71146	8.32%	0.902%
49	11.384	1684	1689	1691	rBV	44984	67269	7.87%	0.853%
50	11.451	1696	1700	1706	rVB	19113	24582	2.88%	0.312%
51	11.616	1722	1727	1736	rBV	100516	124295	14.54%	1.576%
52	11.756	1744	1750	1755	rBV	199720	242821	28.41%	3.078%
53	12.024	1789	1794	1800	rBV	293246	367065	42.94%	4.653%
54	12.091	1800	1805	1809	rBV	28208	34729	4.06%	0.440%
55	12.244	1825	1830	1838	rBV2	368207	469364	54.91%	5.950%
56	12.317	1838	1842	1850	rVB2	35999	53631	6.27%	0.680%
57	12.408	1852	1857	1862	rBV2	17602	23800	2.78%	0.302%
58	12.463	1862	1866	1873	rVB	20032	26587	3.11%	0.337%
59	12.530	1873	1877	1880	rBV	24541	33281	3.89%	0.422%
60	12.555	1880	1881	1886	rVB	15556	13629	1.59%	0.173%
61	12.616	1886	1891	1894	rBV	68761	80533	9.42%	1.021%
62	12.646	1894	1896	1900	rVB	20060	21231	2.48%	0.269%
63	12.701	1900	1905	1913	rBV2	75056	109903	12.86%	1.393%
64	12.920	1933	1941	1945	rBV	43648	59108	6.92%	0.749%
65	12.963	1945	1948	1954	rVV	36955	44566	5.21%	0.565%
66	13.024	1954	1958	1968	rVB3	10095	18988	2.22%	0.241%
67	13.170	1978	1982	1986	rVB	54924	64976	7.60%	0.824%
68	13.256	1992	1996	1998	rBV2	6537	8945	1.05%	0.113%
69	13.292	1998	2002	2007	rVV2	81715	106692	12.48%	1.352%
70	13.426	2020	2024	2029	rVB	19648	25354	2.97%	0.321%
71	13.506	2032	2037	2040	rVV2	9113	12353	1.45%	0.157%
72	13.548	2040	2044	2047	rVV	18908	28281	3.31%	0.359%
73	13.585	2047	2050	2056	rVV4	18303	33822	3.96%	0.429%
74	13.640	2056	2059	2061	rVV2	10072	15511	1.81%	0.197%
75	13.664	2061	2063	2069	rVB2	20231	23662	2.77%	0.300%

LSC Area Percent Report

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Acq On : 06 Dec 2021 20:49
Operator : JC/MD
Sample : M4940-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6

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\82X111721W.M
Title : SW846 8260

76	13.780	2075	2082	2091	rBV	77975	102133	11.95%	1.295%
77	13.926	2100	2106	2111	rBV3	10056	14674	1.72%	0.186%
78	14.079	2126	2131	2136	rBV	14116	17012	1.99%	0.216%
79	14.213	2149	2153	2163	rBV	11279	18983	2.22%	0.241%
80	14.371	2175	2179	2185	rVB	9143	12184	1.43%	0.154%
81	14.640	2219	2223	2228	rVB	7106	8773	1.03%	0.111%
82	14.780	2242	2246	2251	rBV2	9311	12205	1.43%	0.155%

Sum of corrected areas: 7888504

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120621\
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Acq On : 06 Dec 2021 20:49
Operator : JC/MD
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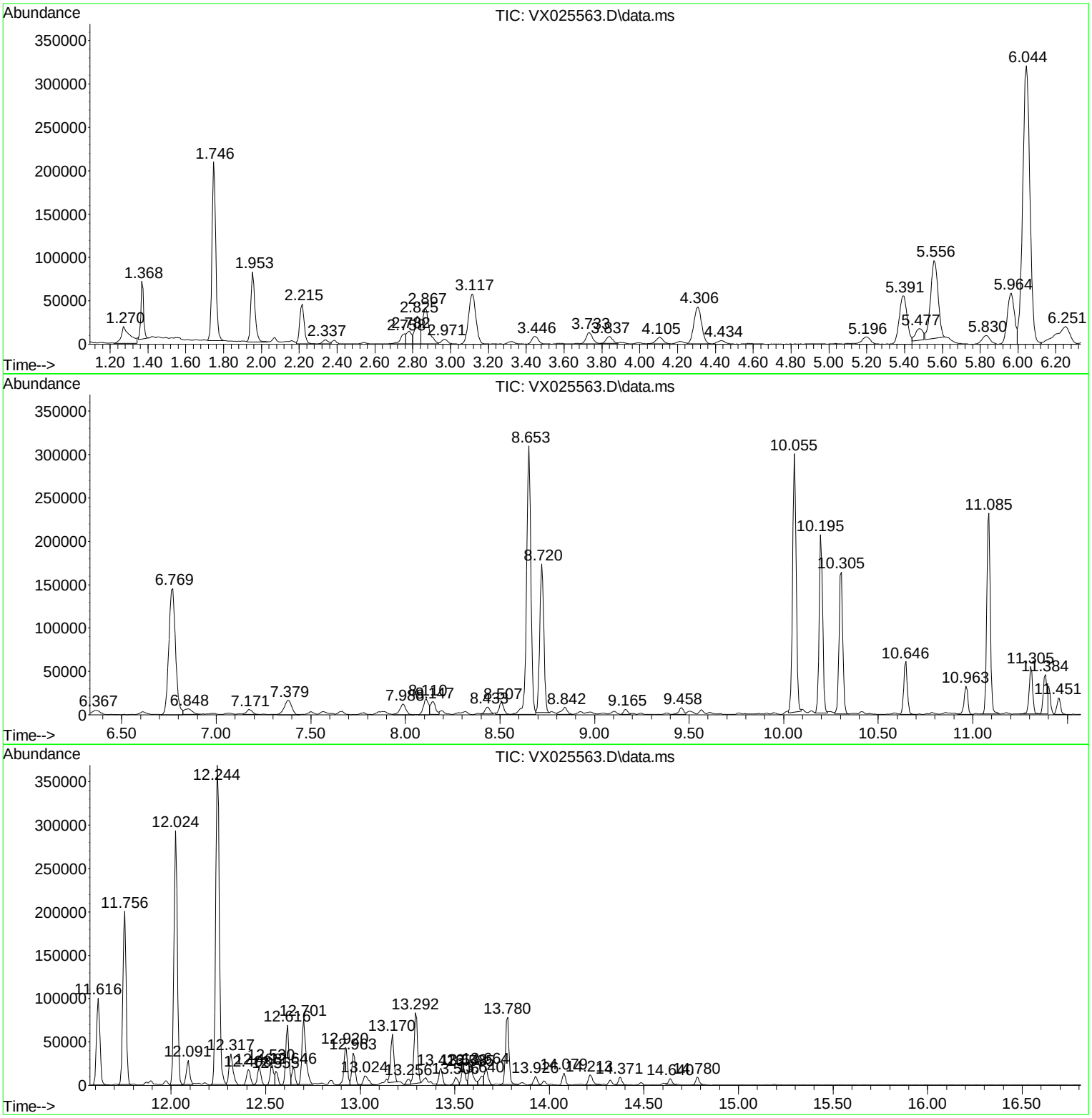
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X111721W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120621\
Data File : VX025563.D
Acq On : 06 Dec 2021 20:49
Operator : JC/MD
Sample : M4940-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X111721W.M
Quant Title : SW846 8260

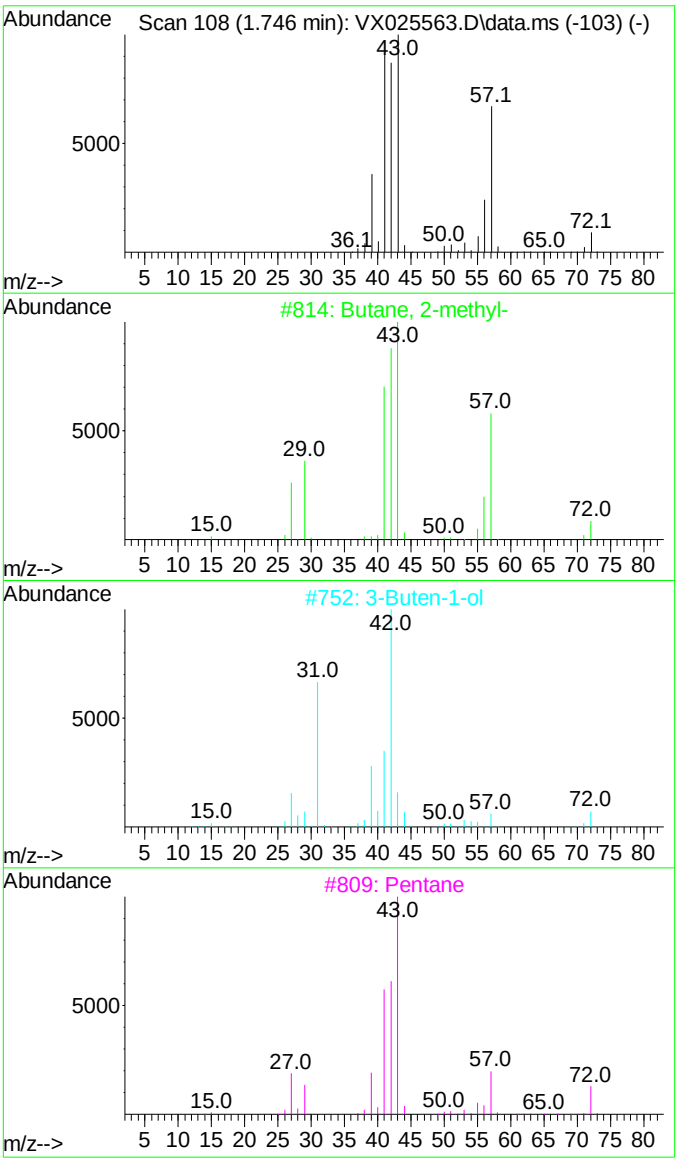
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

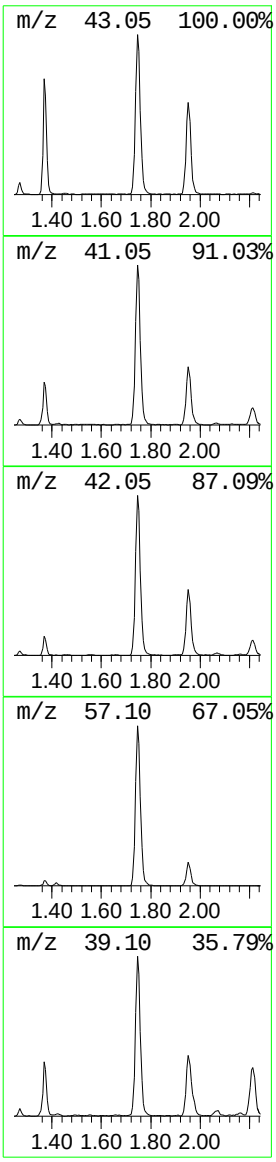
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.746	57.46 ug/l	285186	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	86
2	3-Buten-1-ol	72	C4H8O	000627-27-0	47
3	Pentane	72	C5H12	000109-66-0	33
4	2(3H)-Furanone, dihydro-4-methyl-	100	C5H8O2	001679-49-8	28
5	Butane, 1-chloro-2-methyl-	106	C5H11Cl	000616-13-7	16





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120621\
Data File : VX025563.D
Acq On : 06 Dec 2021 20:49
Operator : JC/MD
Sample : M4940-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X111721W.M
Quant Title : SW846 8260

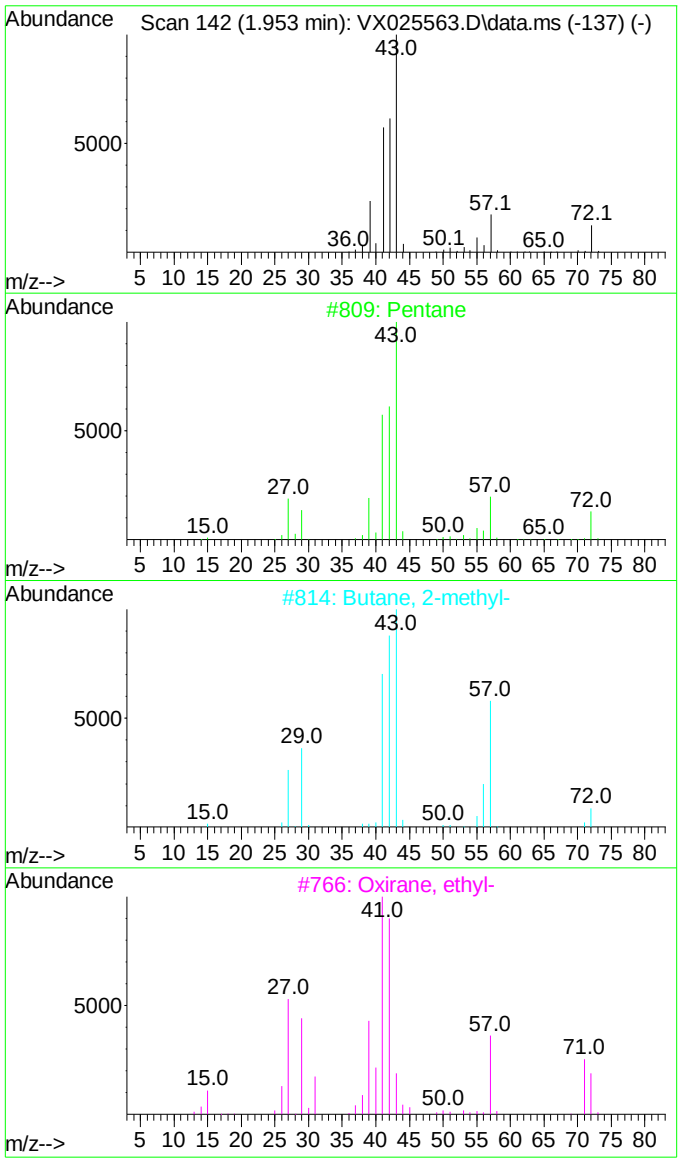
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

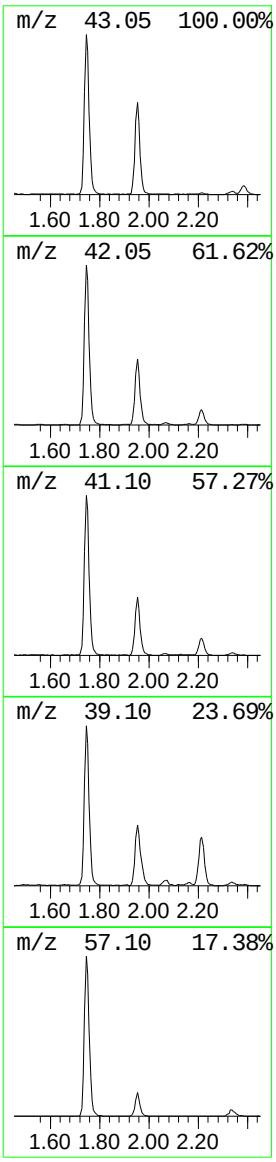
****

Peak Number 2 Pentane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.953	24.57 ug/l	121946	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane	72	C5H12	000109-66-0	91
2	Butane, 2-methyl-	72	C5H12	000078-78-4	64
3	Oxirane, ethyl-	72	C4H8O	000106-88-7	47
4	Isobutyl nitrite	103	C4H9NO2	000542-56-3	17
5	1-Propene, 2-methoxy-	72	C4H8O	000116-11-0	9





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120621\
Data File : VX025563.D
Acq On : 06 Dec 2021 20:49
Operator : JC/MD
Sample : M4940-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6

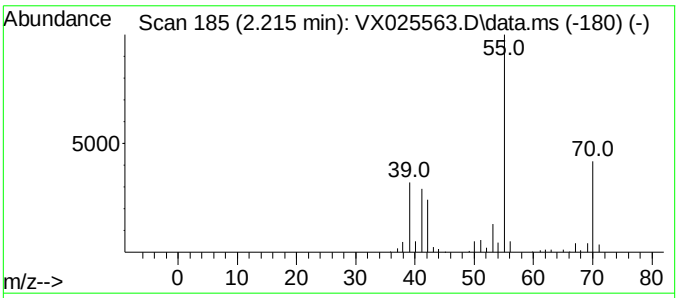
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X111721W.M
Quant Title : SW846 8260

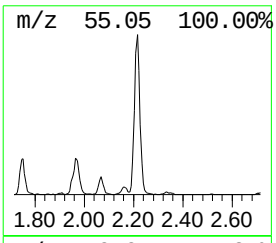
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

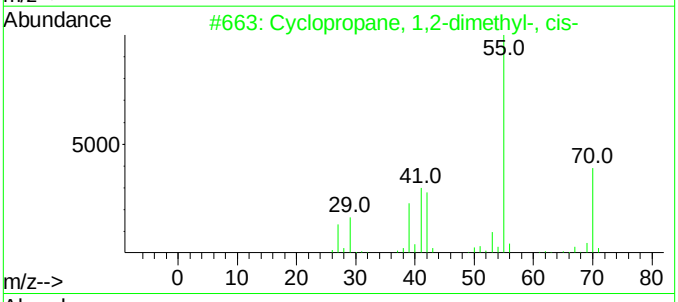
****
Peak Number 3 Cyclopropane, 1,2-dimethyl-... Concentration Rank 9

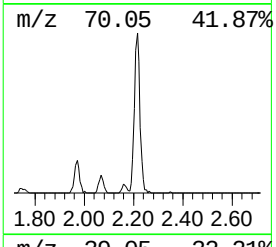
R.T.	EstConc	Area	Relative to ISTD	R.T.
2.215	14.89 ug/l	73888	Pentafluorobenzene	5.556

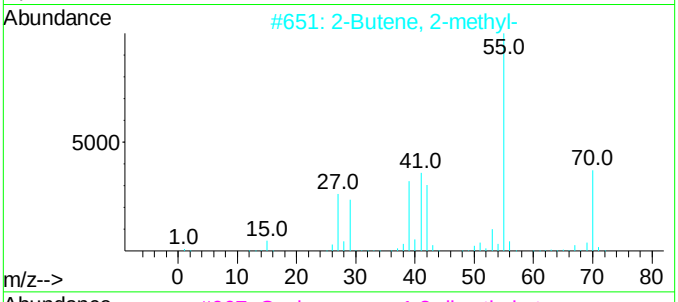
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	91
2	2-Butene, 2-methyl-	70	C5H10	000513-35-9	90
3	Cyclopropane, 1,2-dimethyl-, trans-	70	C5H10	002402-06-4	87
4	2-Pentene, (E)-	70	C5H10	000646-04-8	87
5	1-Butene, 3-methyl-	70	C5H10	000563-45-1	86

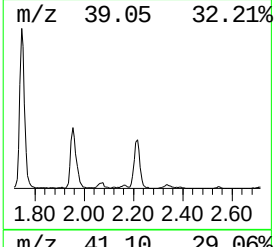


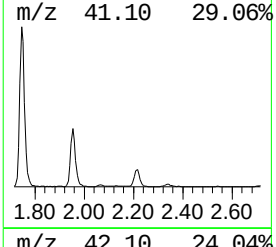


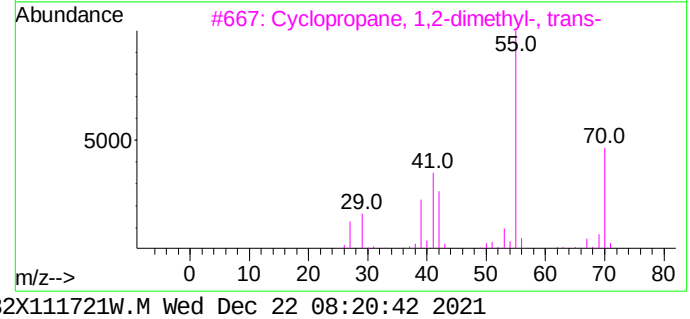


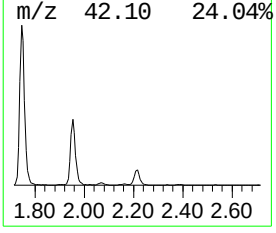












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Data File : VX025563.D
Acq On : 06 Dec 2021 20:49
Operator : JC/MD
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X111721W.M
Quant Title : SW846 8260

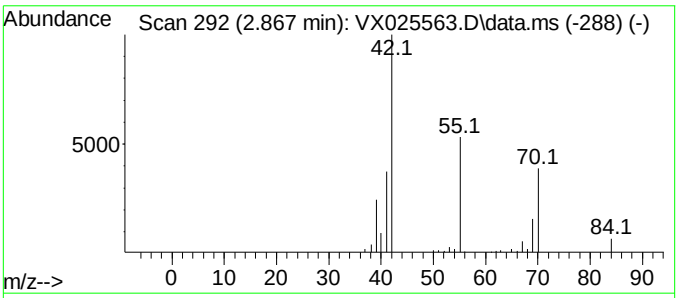
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

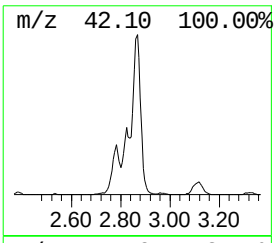
****

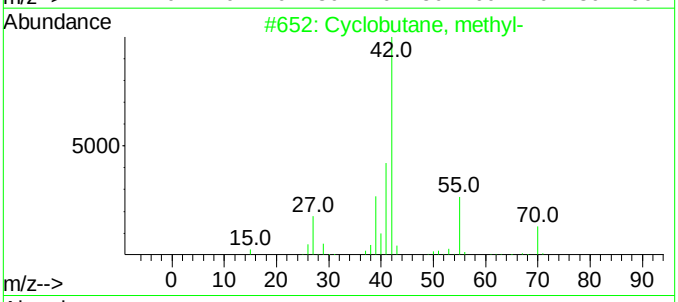
Peak Number 4 Cyclobutane, methyl- Concentration Rank 5

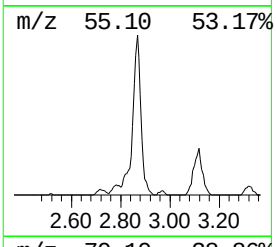
R.T.	EstConc	Area	Relative to ISTD	R.T.
2.867	19.33 ug/l	95964	Pentafluorobenzene	5.556

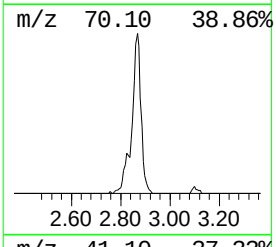
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclobutane, methyl-	70	C5H10	000598-61-8	80
2	Cyclopentane	70	C5H10	000287-92-3	80
3	1-Pentene	70	C5H10	000109-67-1	50
4	3-Buten-1-ol	72	C4H8O	000627-27-0	25
5	2-Pentene, (E)-	70	C5H10	000646-04-8	25

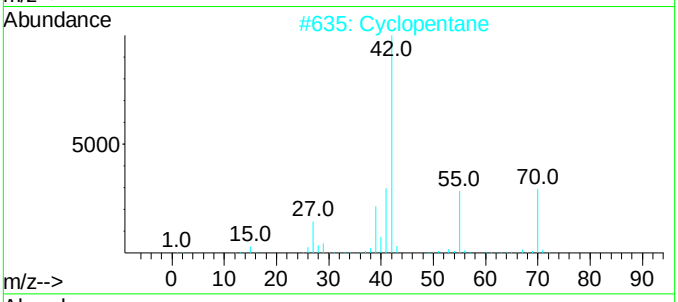


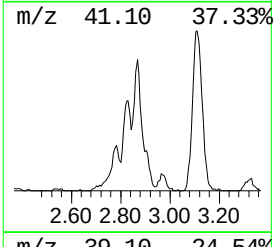


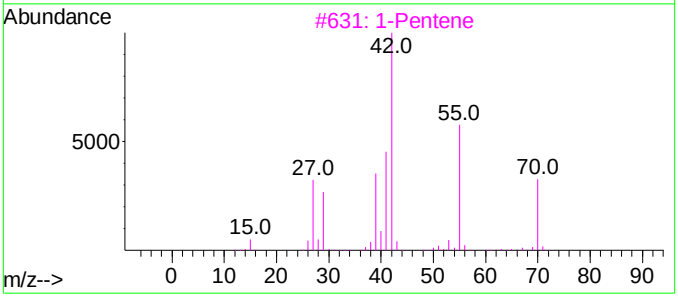


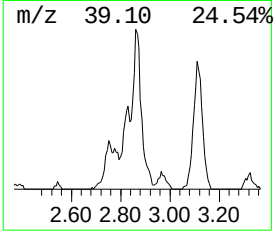












Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120621\
Data File : VX025563.D
Acq On : 06 Dec 2021 20:49
Operator : JC/MD
Sample : M4940-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X111721W.M
Quant Title : SW846 8260

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

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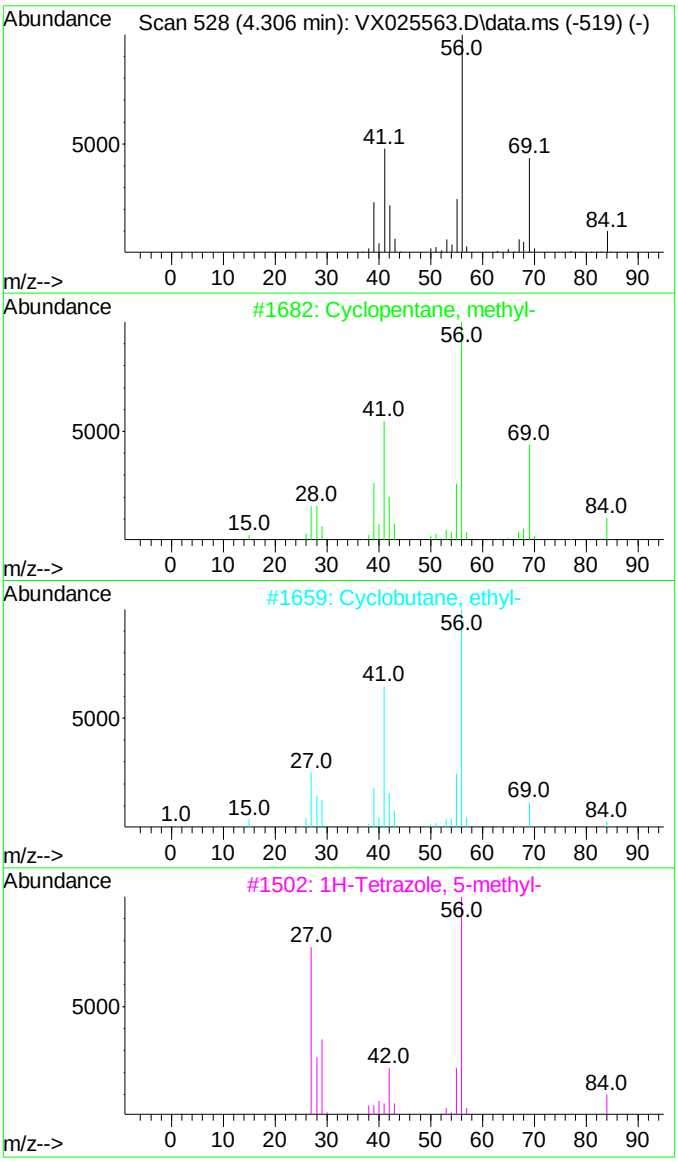
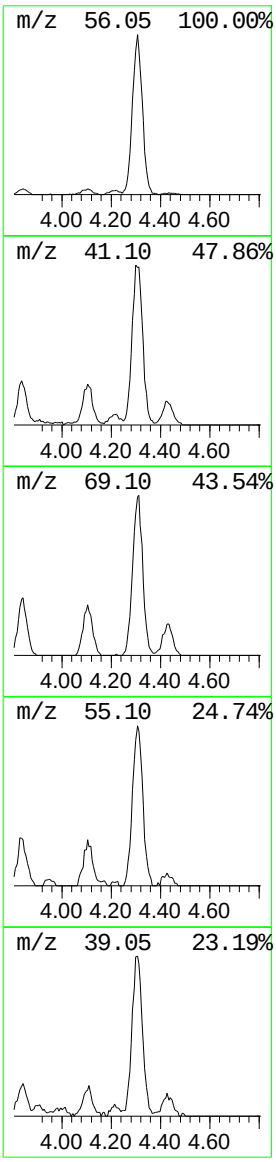
Peak Number 5 Cyclobutane, ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.306	24.70 ug/l	122619	Pentafluorobenzene	5.556

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	72
3	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	50
4	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	47
5	Cyclohexane	84	C6H12	000110-82-7	47





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120621\
Data File : VX025563.D
Acq On : 06 Dec 2021 20:49
Operator : JC/MD
Sample : M4940-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X111721W.M
Quant Title : SW846 8260

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

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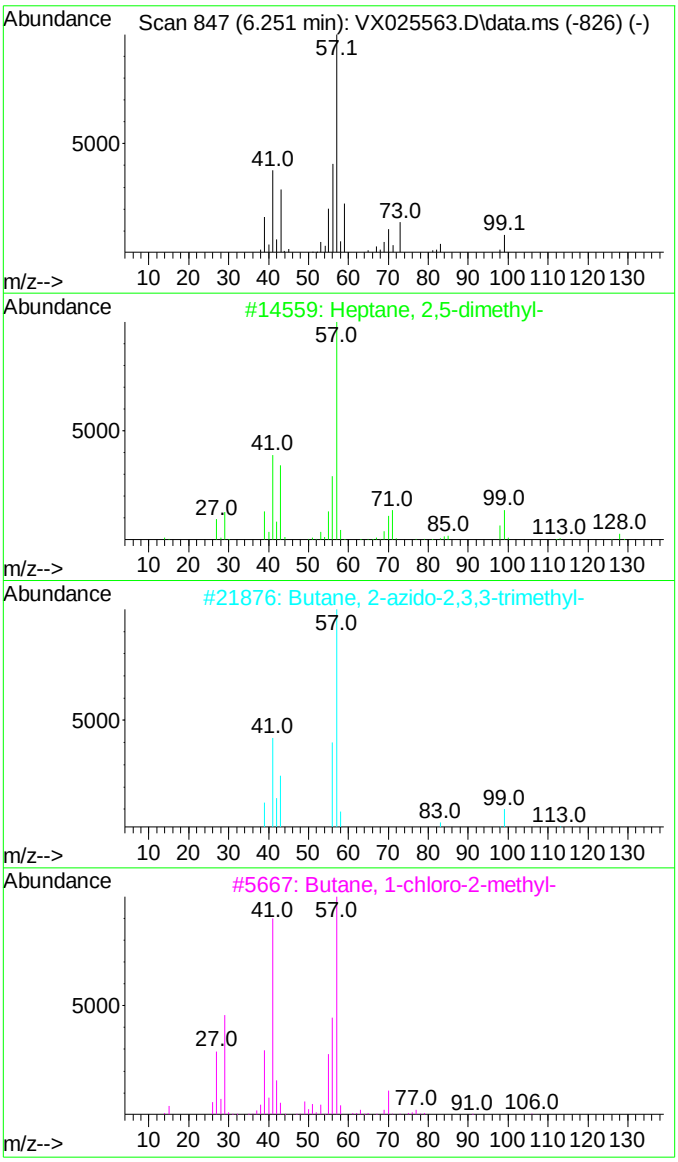
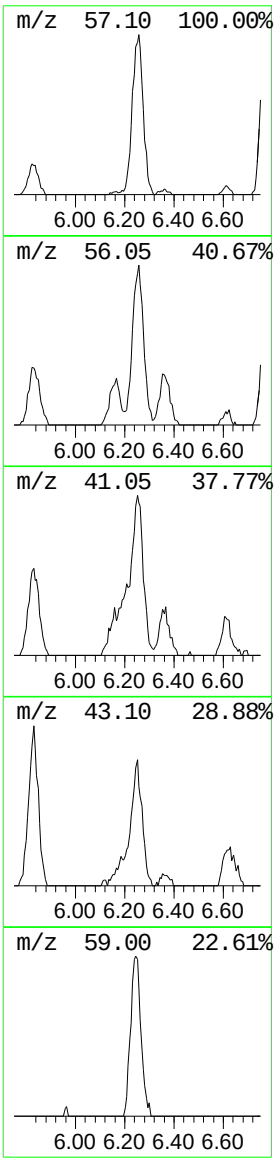
Peak Number 6 Heptane, 2,5-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.251	15.04 ug/l	105898	1,4-Difluorobenzene	6.769

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	53
2	Butane, 2-azido-2,3,3-trimethyl-	141	C7H15N3	051677-41-9	38
3	Butane, 1-chloro-2-methyl-	106	C5H11Cl	000616-13-7	38
4	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	37
5	Hexane, 2,2,5,5-tetramethyl-	142	C10H22	001071-81-4	37





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120621\
Data File : VX025563.D
Acq On : 06 Dec 2021 20:49
Operator : JC/MD
Sample : M4940-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X111721W.M
Quant Title : SW846 8260

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

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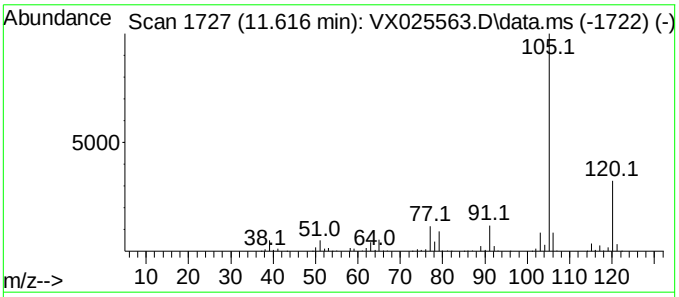
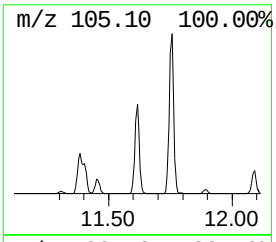
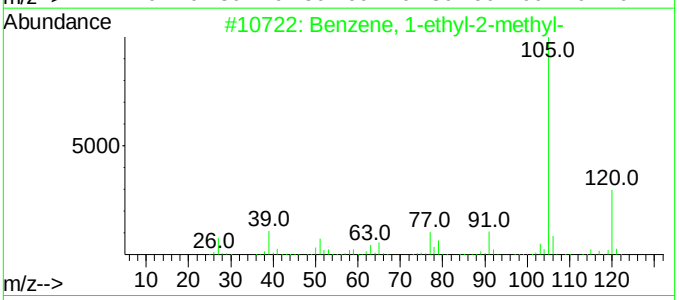
Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 6

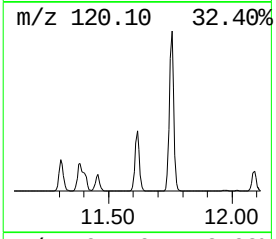
R.T.	EstConc	Area	Relative to ISTD	R.T.
11.616	16.93 ug/l	124295	1,4-Dichlorobenzene-d4	12.024

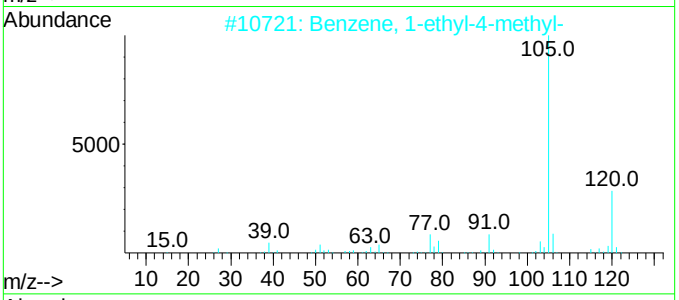
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5	Mesitylene	120	C9H12	000108-67-8	91

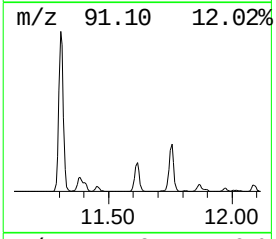


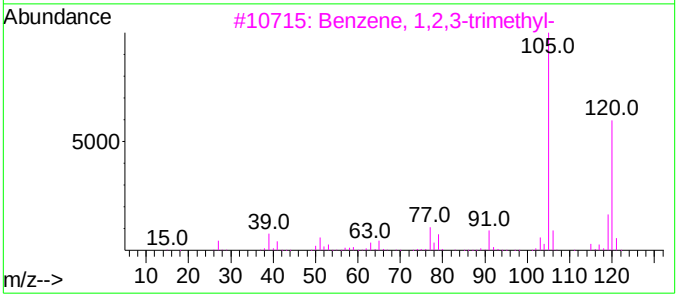


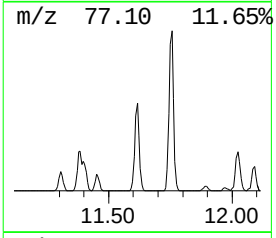


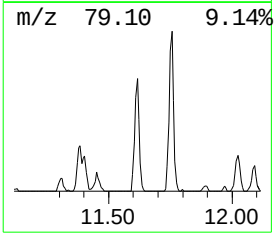












Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120621\
Data File : VX025563.D
Acq On : 06 Dec 2021 20:49
Operator : JC/MD
Sample : M4940-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X111721W.M
Quant Title : SW846 8260

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

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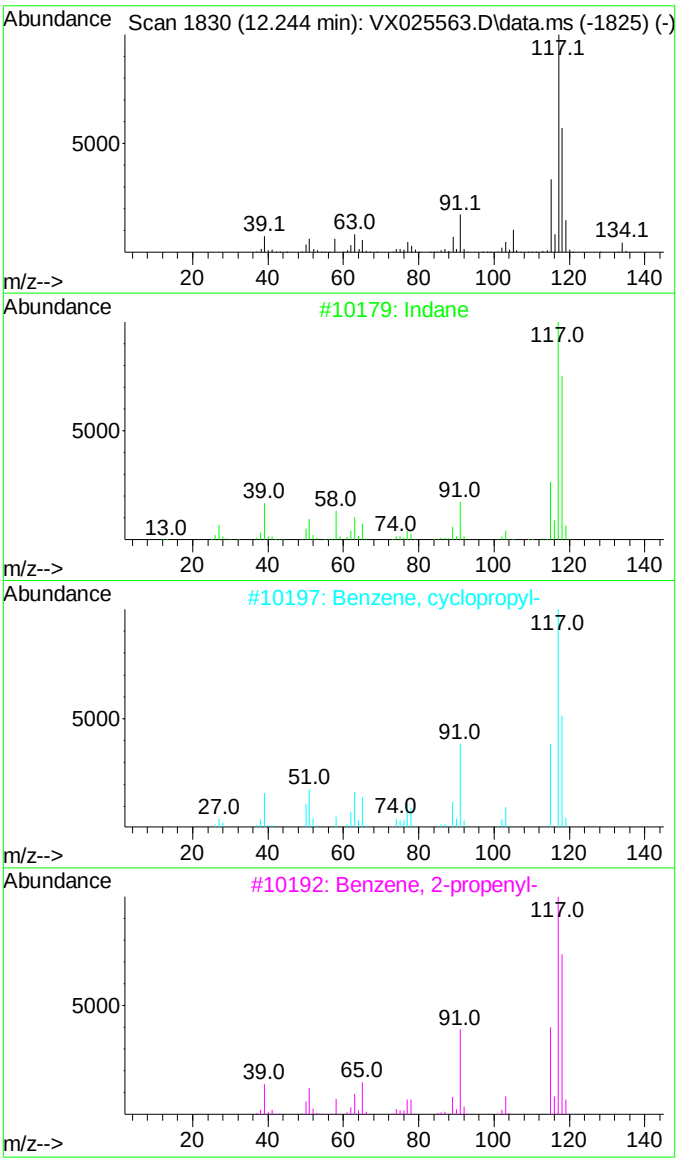
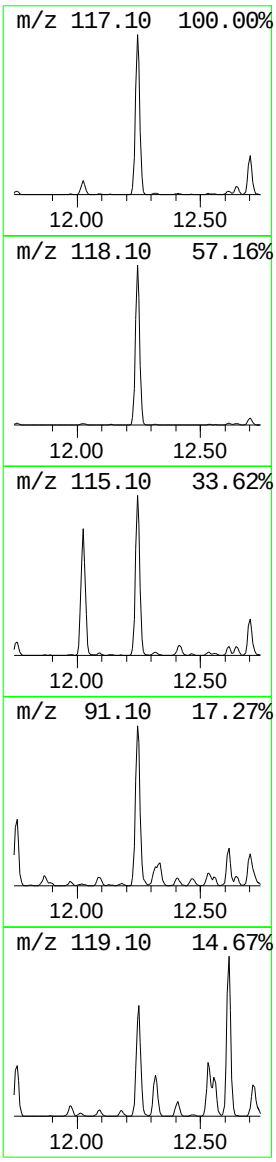
Peak Number 8 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	63.93 ug/l	469364	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	76
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
4	Deltacyclene	118	C9H10	007785-10-6	58
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	55





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120621\
Data File : VX025563.D
Acq On : 06 Dec 2021 20:49
Operator : JC/MD
Sample : M4940-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X111721W.M
Quant Title : SW846 8260

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

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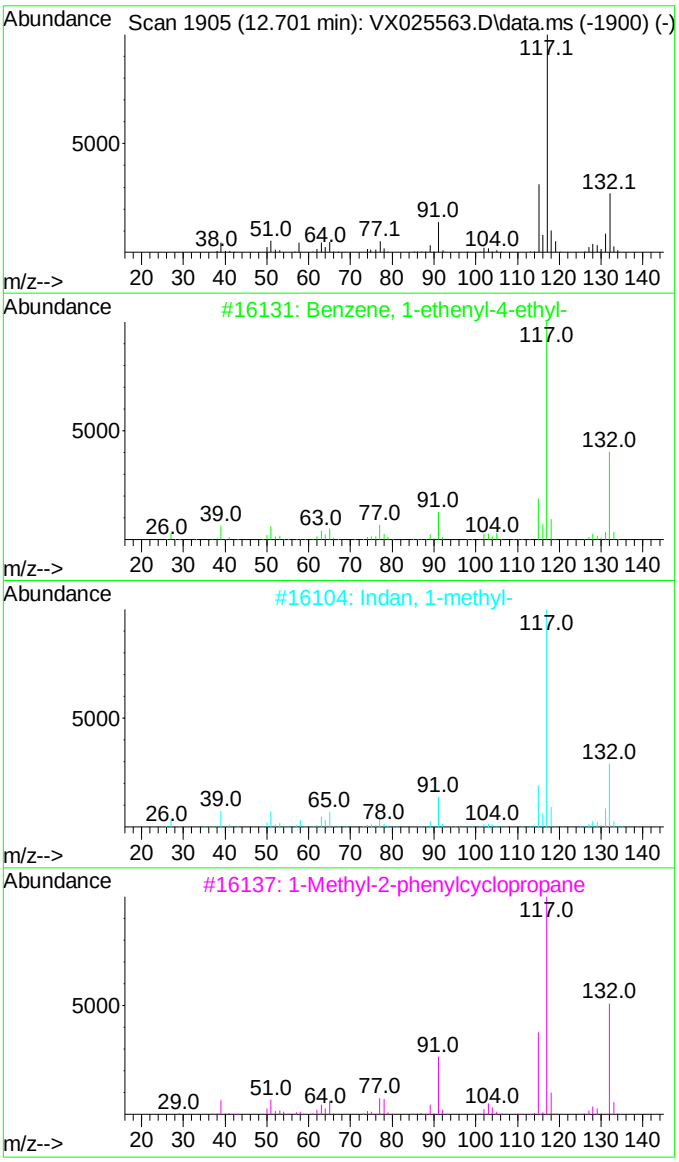
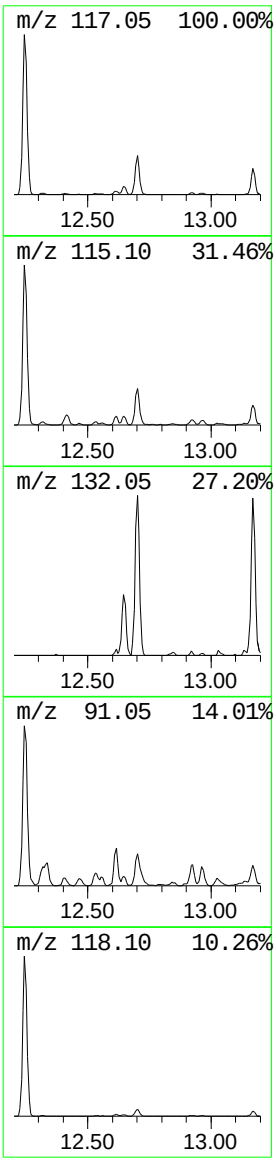
Peak Number 9 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	14.97 ug/l	109903	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	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	86
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	83





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120621\
Data File : VX025563.D
Acq On : 06 Dec 2021 20:49
Operator : JC/MD
Sample : M4940-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X111721W.M
Quant Title : SW846 8260

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

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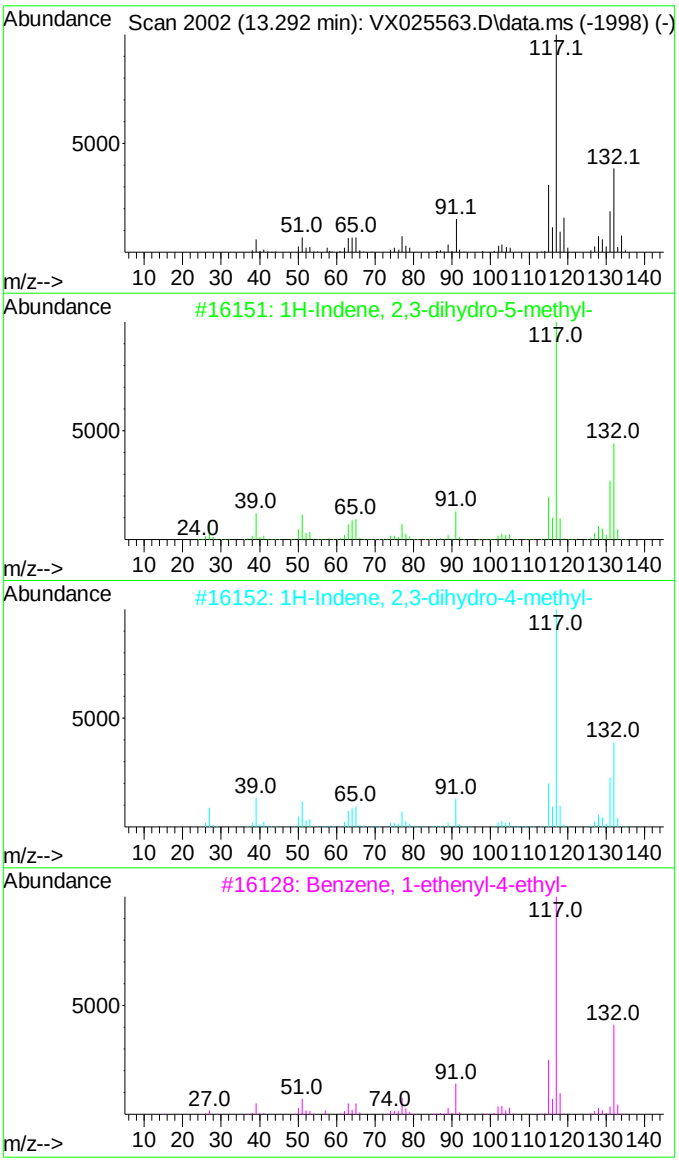
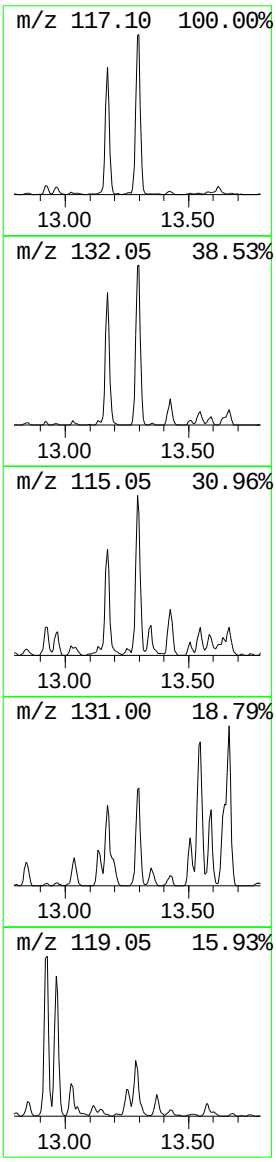
Peak Number 10 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	14.53 ug/l	106692	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	87
2	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
3	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
4	1-Phenyl-1-butene	132	C10H12	000824-90-8	86
5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	86





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120621\
Data File : VX025563.D
Acq On : 06 Dec 2021 20:49
Operator : JC/MD
Sample : M4940-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X111721W.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-	1.746	57.5	ug/l	285186	1	5.556	248167	50.0
Pentane	1.953	24.6	ug/l	121946	1	5.556	248167	50.0
Cyclopropane, 1...	2.215	14.9	ug/l	73888	1	5.556	248167	50.0
Cyclobutane, me...	2.867	19.3	ug/l	95964	1	5.556	248167	50.0
Cyclobutane, et...	4.306	24.7	ug/l	122619	1	5.556	248167	50.0
Heptane, 2,5-di...	6.251	15.0	ug/l	105898	2	6.769	352088	50.0
Benzene, 1-ethy...	11.616	16.9	ug/l	124295	4	12.024	367065	50.0
Indane	12.244	63.9	ug/l	469364	4	12.024	367065	50.0
Benzene, 1-ethe...	12.701	15.0	ug/l	109903	4	12.024	367065	50.0
1H-Indene, 2,3-...	13.292	14.5	ug/l	106692	4	12.024	367065	50.0