

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100623\
 Data File : VX038123.D
 Acq On : 06 Oct 2023 12:14
 Operator : JC/MD
 Sample : 04723-02 20X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-5B

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\82X100523W.M
 Title : SW846 8260

Signal : TIC: VX038123.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.373	43	47	53	rBV	40158	41838	5.12%	0.390%
2	1.453	56	60	69	rVB	50794	57421	7.03%	0.535%
3	1.562	69	78	83	rBV2	15041	37725	4.62%	0.352%
4	1.745	103	108	121	rVB	243304	344421	42.16%	3.209%
5	1.952	137	142	152	rVB2	57175	97880	11.98%	0.912%
6	2.068	156	161	168	rBV	18399	28139	3.44%	0.262%
7	2.166	173	177	180	rVV	12309	20780	2.54%	0.194%
8	2.215	180	185	193	rVB	51218	73441	8.99%	0.684%
9	2.343	198	206	218	rVB	12962	27724	3.39%	0.258%
10	2.782	269	278	281	rBV	39609	84978	10.40%	0.792%
11	2.824	281	285	289	rVV	46842	99194	12.14%	0.924%
12	2.867	289	292	302	rVB2	34770	66545	8.15%	0.620%
13	3.099	322	330	346	rBV	61449	139453	17.07%	1.299%
14	3.318	360	366	375	rVB4	5268	12316	1.51%	0.115%
15	3.446	380	387	396	rVB2	5811	13241	1.62%	0.123%
16	3.733	427	434	444	rVB	10873	25104	3.07%	0.234%
17	3.842	444	452	458	rBV3	10209	26487	3.24%	0.247%
18	3.897	458	461	468	rVB2	3644	8423	1.03%	0.078%
19	4.104	482	495	504	rVB5	8540	24792	3.03%	0.231%
20	4.208	504	512	518	rVV3	4331	11597	1.42%	0.108%
21	4.306	518	528	539	rVV	68622	194777	23.84%	1.815%
22	4.428	539	548	561	rVB4	6439	21102	2.58%	0.197%
23	4.562	561	570	586	rBV2	6474	23761	2.91%	0.221%
24	5.135	644	664	665	rBV4	2580	8396	1.03%	0.078%
25	5.196	665	674	689	rVB2	12931	42852	5.25%	0.399%
26	5.385	694	705	714	rBV	75109	224425	27.47%	2.091%
27	5.476	714	720	725	rVV2	17200	39157	4.79%	0.365%
28	5.555	725	733	742	rVV	106044	281485	34.46%	2.623%
29	5.830	766	778	790	rBV5	12443	40540	4.96%	0.378%
30	5.958	790	799	812	rBV	84350	228629	27.99%	2.130%
31	6.165	824	833	843	rBV6	12894	50040	6.13%	0.466%
32	6.257	843	848	856	rVV4	13083	34427	4.21%	0.321%
33	6.354	858	864	876	rVB4	11656	33775	4.13%	0.315%
34	6.763	921	931	942	rBV	202179	501405	61.38%	4.672%
35	7.171	988	998	1006	rBV4	11572	27417	3.36%	0.255%
36	7.378	1021	1032	1043	rBV5	32642	92322	11.30%	0.860%

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37	7.659	1072	1078	1086	rVB2	7736	14991	1.84%	0.140%
38	7.775	1090	1097	1103	rBV4	4092	8237	1.01%	0.077%
39	8.146	1153	1158	1163	rVV	23641	41902	5.13%	0.390%
40	8.311	1180	1185	1193	rVB5	4135	9054	1.11%	0.084%
41	8.506	1211	1217	1224	rVB3	14138	25778	3.16%	0.240%
42	8.604	1226	1233	1235	rBV3	5673	10170	1.24%	0.095%
43	8.652	1235	1241	1248	rVV	413108	674654	82.59%	6.287%
44	8.720	1248	1252	1259	rVB	58481	91271	11.17%	0.851%
45	8.927	1279	1286	1290	rBV2	5681	9051	1.11%	0.084%
46	9.104	1307	1315	1320	rVV3	5725	10677	1.31%	0.099%
47	9.165	1320	1325	1335	rVB	16903	28894	3.54%	0.269%
48	9.500	1377	1380	1386	rVB4	5369	10292	1.26%	0.096%
49	10.055	1465	1471	1480	rBV	461572	626057	76.64%	5.834%
50	10.195	1489	1494	1503	rVB	331393	422309	51.70%	3.935%
51	10.299	1504	1511	1524	rVB	485362	679880	83.23%	6.335%
52	10.640	1562	1567	1578	rVB	165179	220728	27.02%	2.057%
53	10.963	1614	1620	1625	rBV	33795	41961	5.14%	0.391%
54	11.079	1634	1639	1652	rVB	376594	497397	60.89%	4.635%
55	11.304	1671	1676	1682	rBV	132838	159646	19.54%	1.488%
56	11.402	1682	1692	1697	rVV	107315	194161	23.77%	1.809%
57	11.451	1697	1700	1708	rVB	21986	28038	3.43%	0.261%
58	11.615	1719	1727	1735	rBV	145534	193678	23.71%	1.805%
59	11.749	1744	1749	1757	rBV	657200	816876	100.00%	7.612%
60	11.865	1763	1768	1770	rBV	9866	13012	1.59%	0.121%
61	11.890	1770	1772	1777	rVB	12683	13976	1.71%	0.130%
62	11.969	1781	1785	1788	rBV	9008	10668	1.31%	0.099%
63	12.024	1788	1794	1800	rVV	524769	672455	82.32%	6.266%
64	12.085	1800	1804	1810	rVB	147295	179104	21.93%	1.669%
65	12.243	1823	1830	1837	rBV2	155384	211296	25.87%	1.969%
66	12.310	1837	1841	1851	rVB3	54649	83213	10.19%	0.775%
67	12.408	1851	1857	1862	rBV2	13385	18207	2.23%	0.170%
68	12.463	1862	1866	1871	rVB	40276	48212	5.90%	0.449%
69	12.530	1872	1877	1879	rBV	90734	113041	13.84%	1.053%
70	12.554	1879	1881	1886	rVB	74720	78885	9.66%	0.735%
71	12.609	1886	1890	1894	rBV	134116	171209	20.96%	1.595%
72	12.646	1894	1896	1900	rVB	33556	34153	4.18%	0.318%
73	12.701	1900	1905	1912	rBV2	89226	128467	15.73%	1.197%
74	12.847	1924	1929	1934	rVB2	37362	47350	5.80%	0.441%
75	12.920	1934	1941	1944	rBV	89067	105172	12.87%	0.980%
76	12.963	1944	1948	1954	rVB	120981	142499	17.44%	1.328%

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77	13.024	1954	1958	1967	rVB4	10986	18994	2.33%	0.177%
78	13.133	1971	1976	1978	rBV2	9220	12176	1.49%	0.113%
79	13.170	1978	1982	1992	rVB	81152	110697	13.55%	1.032%
80	13.255	1992	1996	1998	rBV2	10756	16295	1.99%	0.152%
81	13.292	1998	2002	2007	rVB2	144733	187305	22.93%	1.745%
82	13.420	2019	2023	2030	rVB	16409	23195	2.84%	0.216%
83	13.505	2030	2037	2040	rBV3	12634	18139	2.22%	0.169%
84	13.542	2040	2043	2046	rVV	27381	36630	4.48%	0.341%
85	13.585	2046	2050	2055	rVV3	31533	59126	7.24%	0.551%
86	13.639	2055	2059	2060	rVV2	17215	22711	2.78%	0.212%
87	13.664	2060	2063	2069	rVB2	32587	45791	5.61%	0.427%
88	13.773	2074	2081	2090	rBV	54920	74281	9.09%	0.692%
89	13.865	2090	2096	2100	rBV4	5005	9495	1.16%	0.088%
90	13.926	2100	2106	2110	rBV3	10731	16302	2.00%	0.152%
91	13.969	2110	2113	2117	rVB2	9586	11395	1.39%	0.106%
92	14.072	2125	2130	2135	rBV	18249	23779	2.91%	0.222%
93	14.212	2149	2153	2160	rBV2	16146	25470	3.12%	0.237%
94	14.371	2176	2179	2184	rVB2	10728	13123	1.61%	0.122%
95	14.779	2241	2246	2254	rVB	26891	33864	4.15%	0.316%

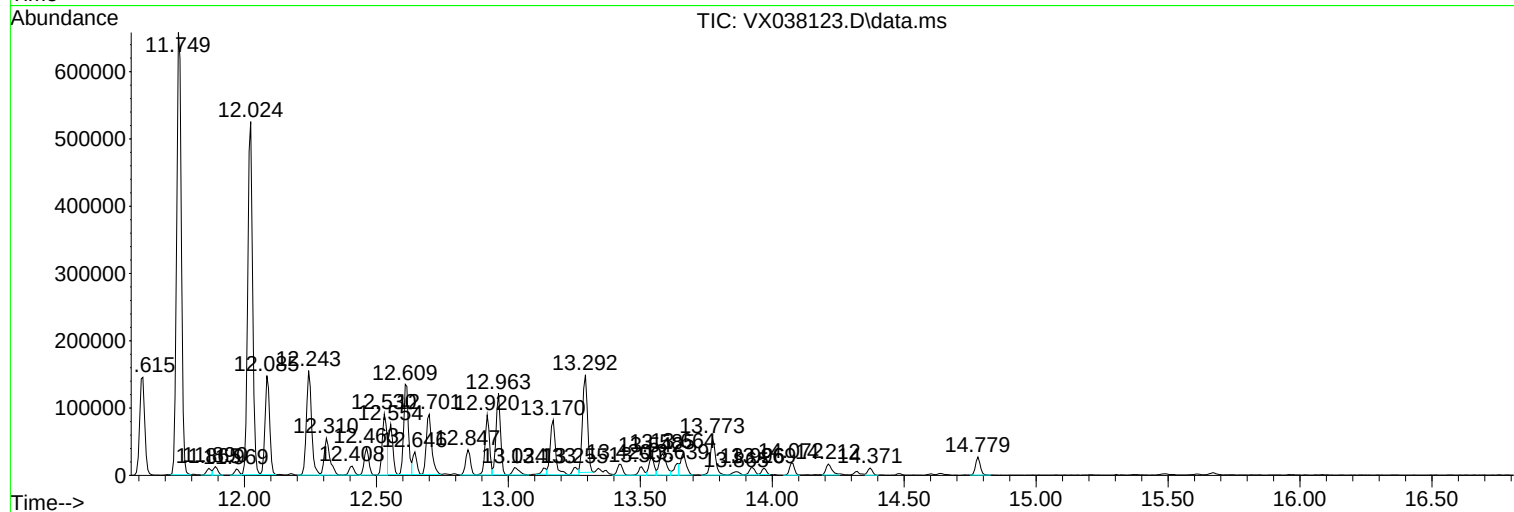
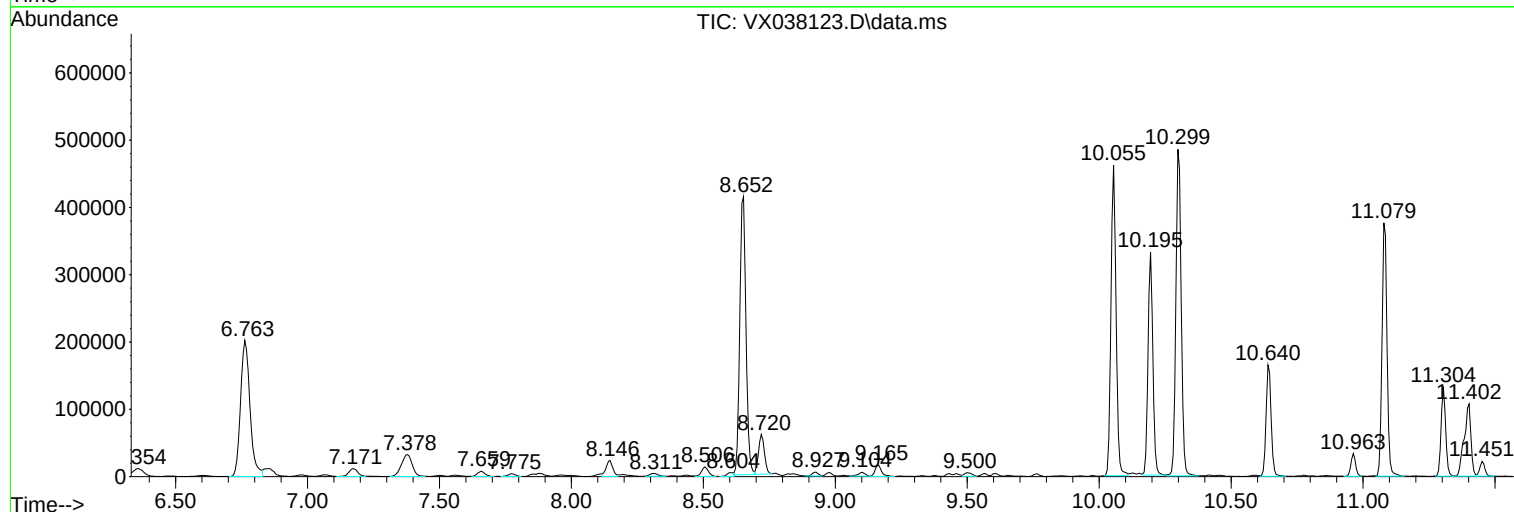
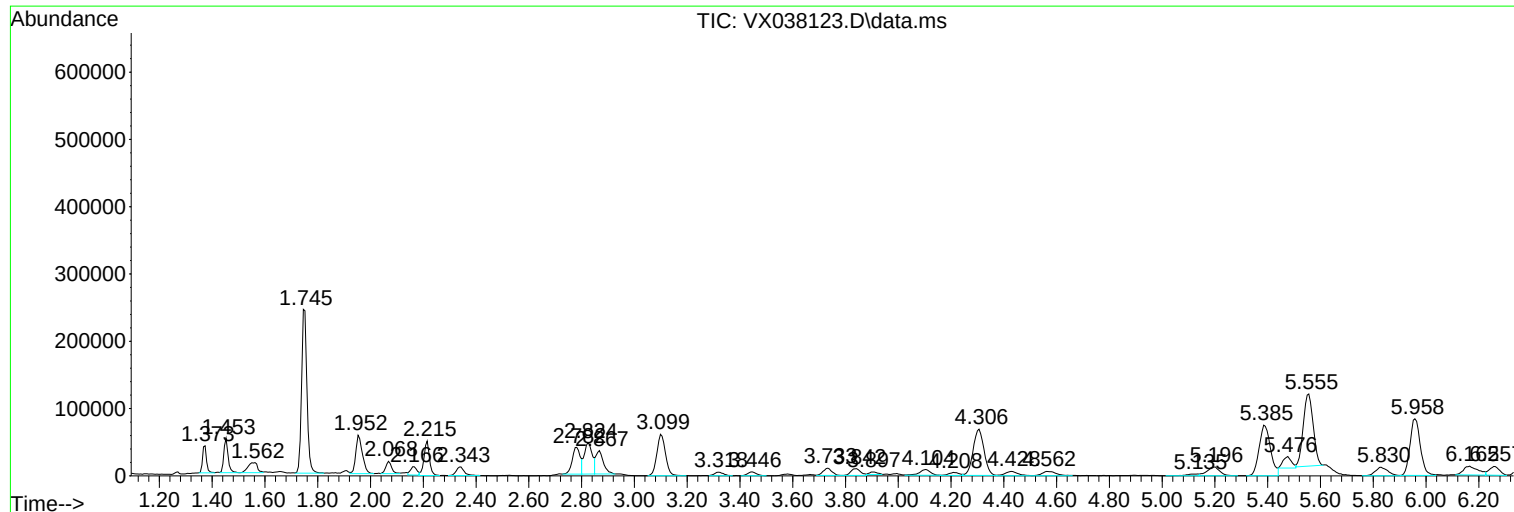
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.M
Quant Title : SW846 8260

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



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Instrument :
MSVOA_X
ClientSampleId :
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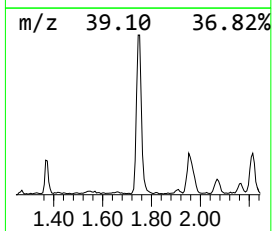
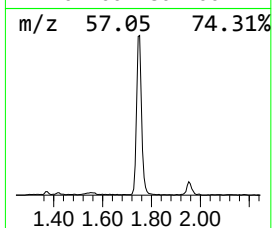
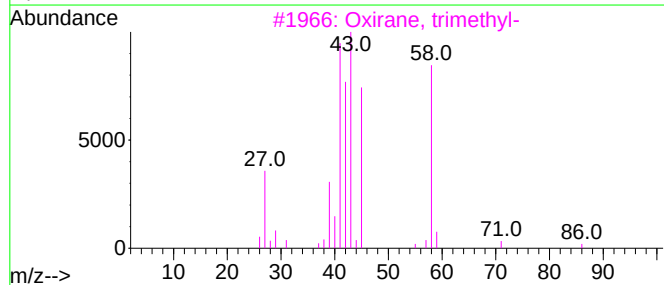
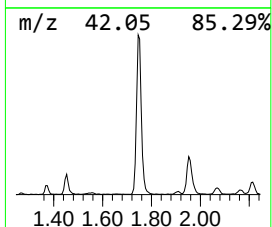
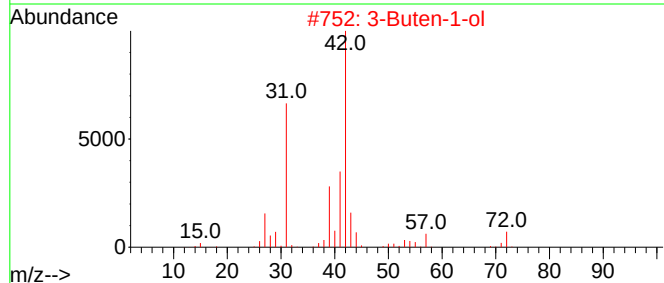
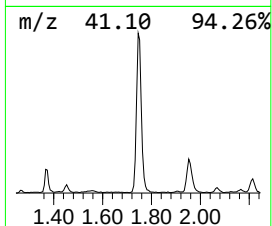
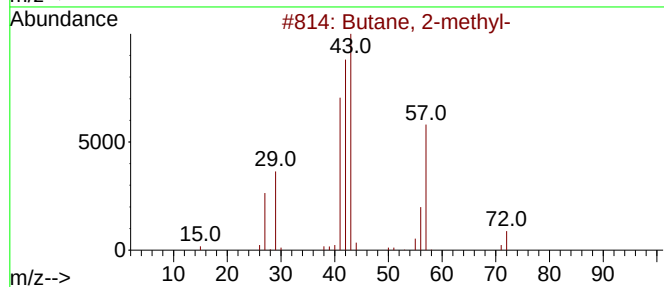
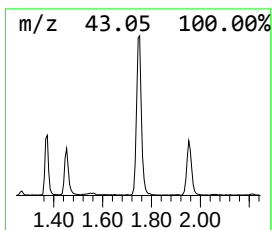
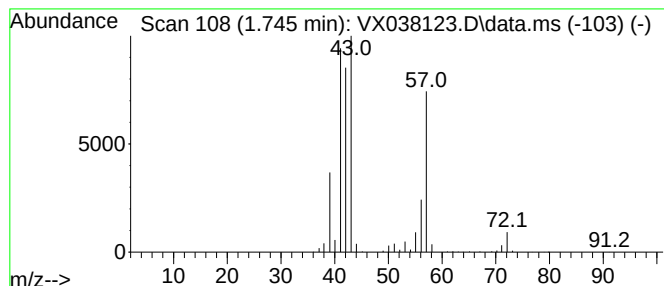
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.745	61.18 ug/l	344421	Pentafluorobenzene	5.555

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	43
3			Oxirane, trimethyl-	86	C5H10O	005076-19-7	40
4			Pentane	72	C5H12	000109-66-0	28
5			2(3H)-Furanone, dihydro-4-methyl-	100	C5H8O2	001679-49-8	25



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Instrument :
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ClientSampleId :
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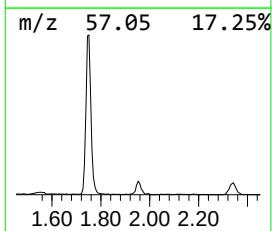
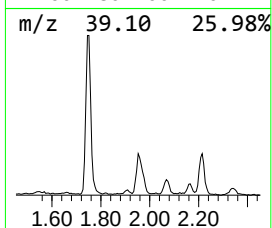
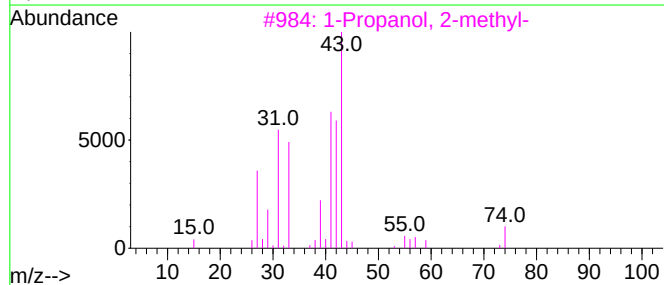
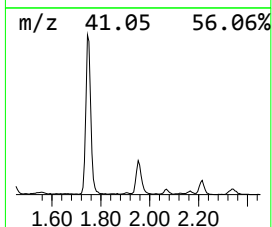
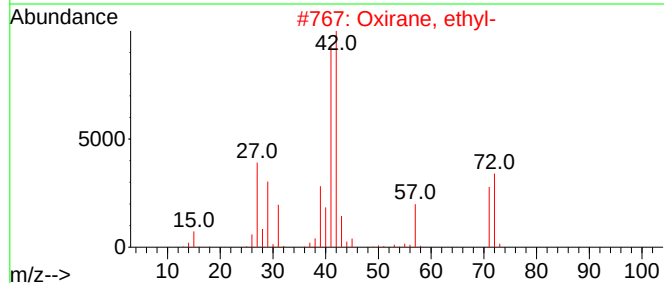
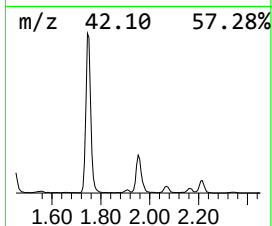
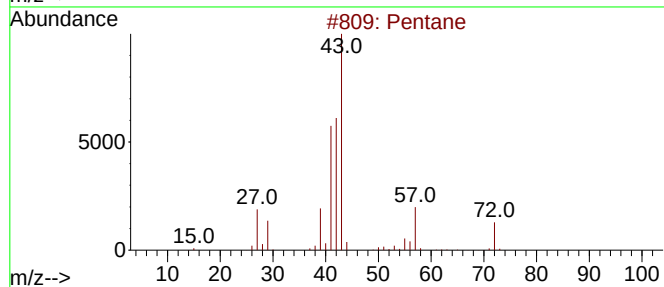
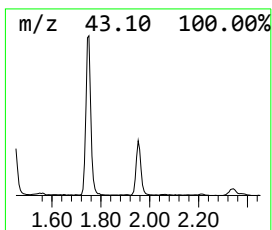
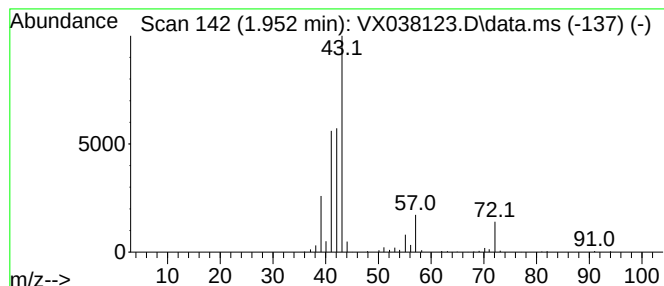
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.M
Quant Title : SW846 8260

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

Peak Number 2 Pentane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.952	17.39 ug/l	97880	Pentafluorobenzene	5.555

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane	72	C5H12	000109-66-0	91
2			Oxirane, ethyl-	72	C4H8O	000106-88-7	50
3			1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	45
4			Butane, 2-methyl-	72	C5H12	000078-78-4	39
5			Tetrahydrofuran	72	C4H8O	000109-99-9	38



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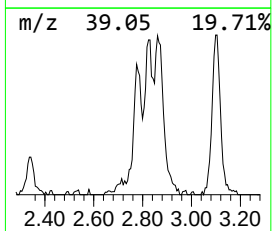
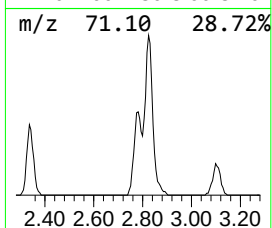
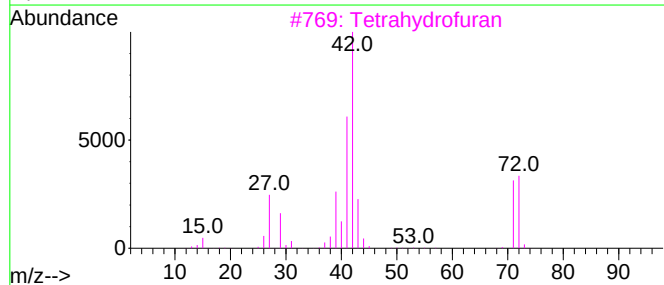
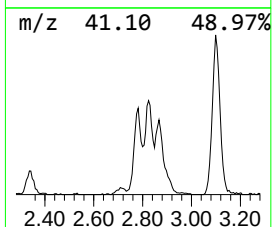
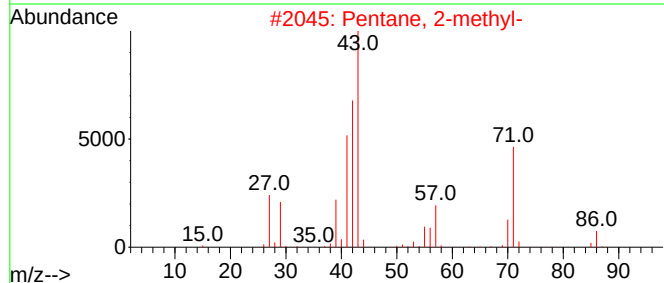
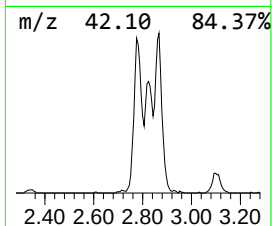
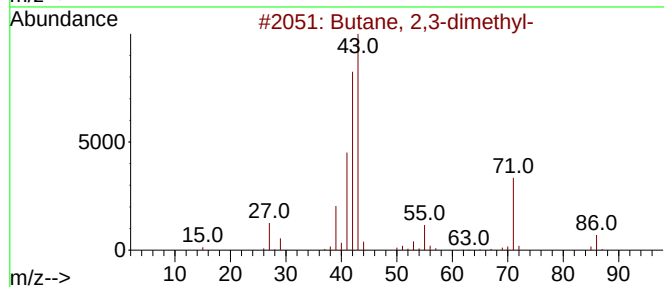
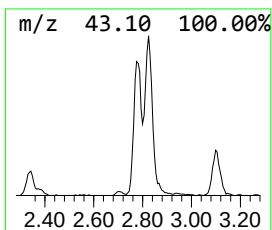
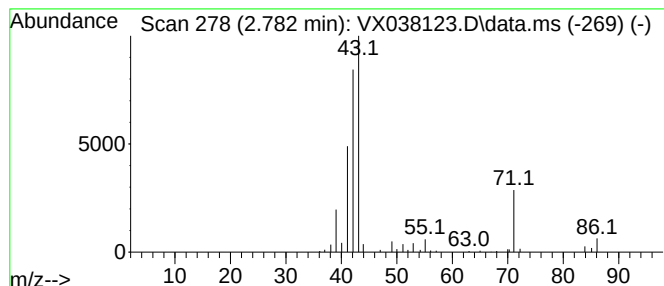
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.M
Quant Title : SW846 8260

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

Peak Number 3 Butane, 2,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.782	15.09 ug/l	84978	Pentafluorobenzene	5.555

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	87
2			Pentane, 2-methyl-	86	C6H14	000107-83-5	40
3			Tetrahydrofuran	72	C4H8O	000109-99-9	36
4			Pentane	72	C5H12	000109-66-0	17
5			Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	10



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Acq On : 06 Oct 2023 12:14
Operator : JC/MD
Sample : 04723-02 20X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-5B

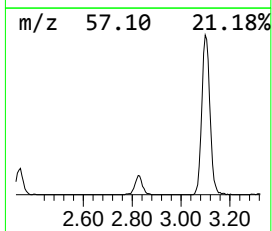
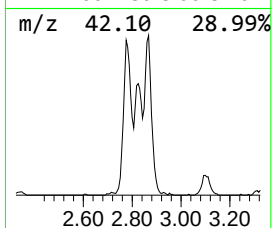
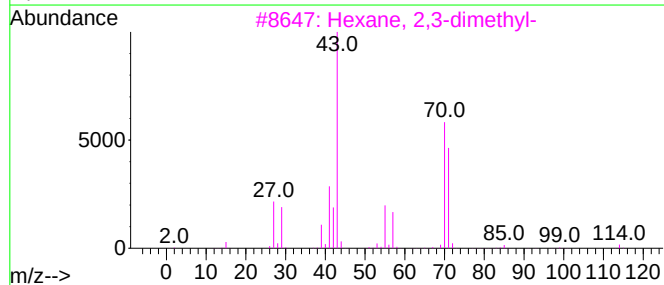
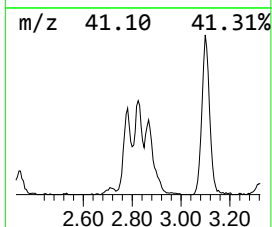
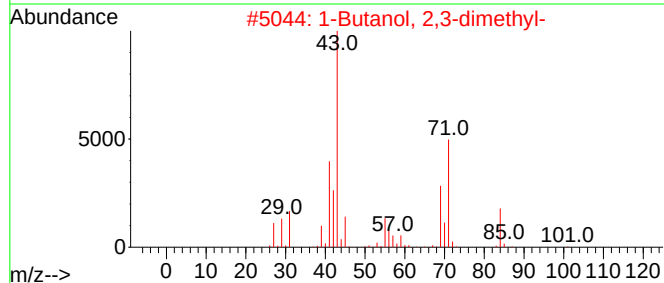
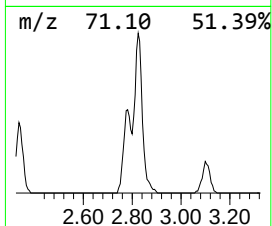
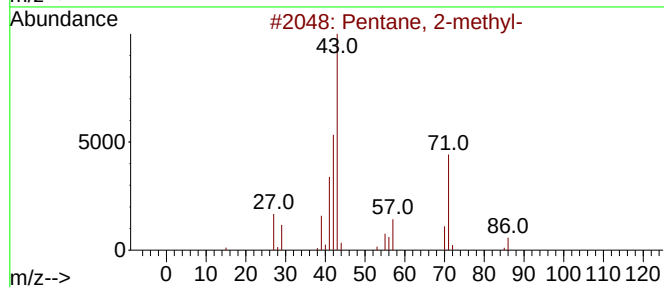
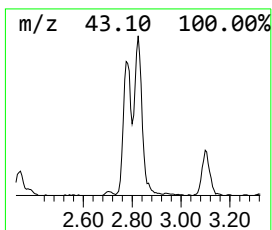
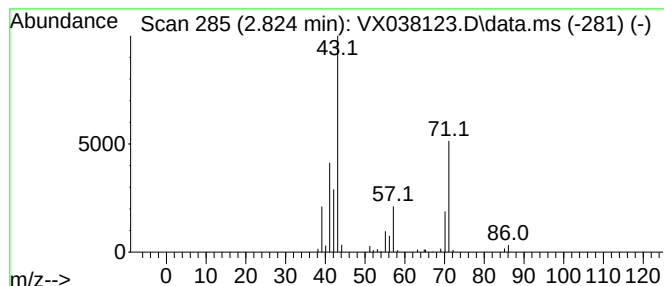
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Quant Title : SW846 8260

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

Peak Number 4 Pentane, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.824	17.62 ug/l	99194	Pentafluorobenzene	5.555

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2-methyl-	86	C6H14	000107-83-5	59
2			1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	40
3			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	38
4			Butane, 1-chloro-3-methyl-	106	C5H11Cl	000107-84-6	37
5			Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	33



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100623\
Data File : VX038123.D
Acq On : 06 Oct 2023 12:14
Operator : JC/MD
Sample : 04723-02 20X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-5B

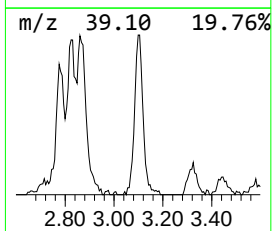
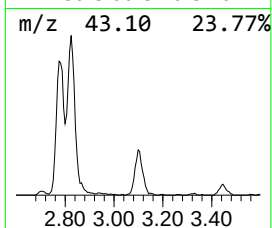
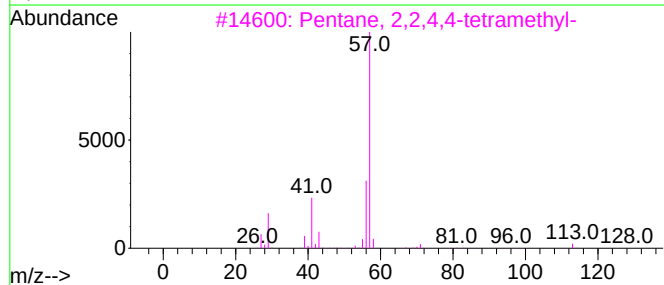
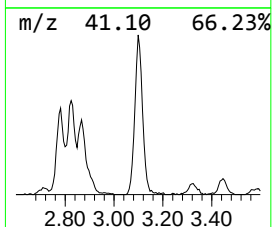
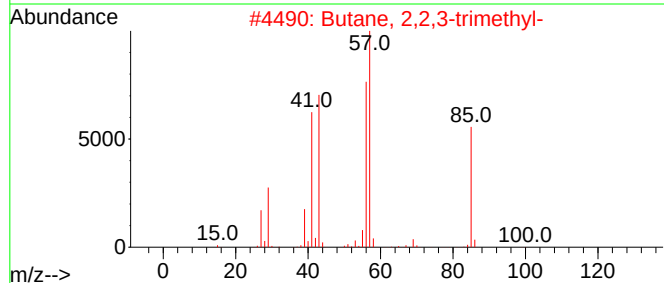
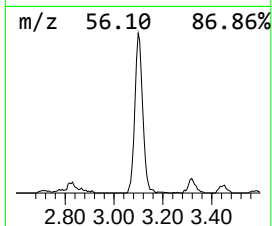
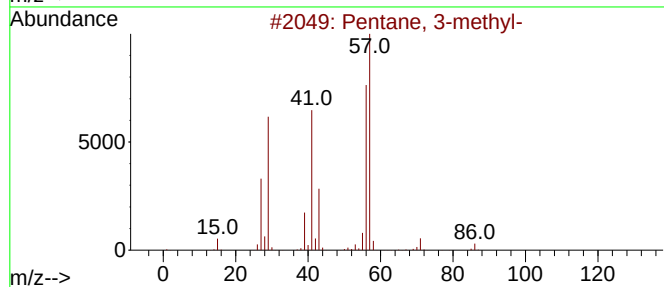
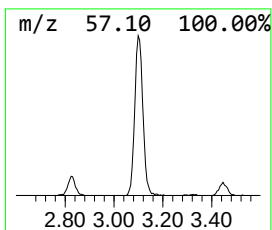
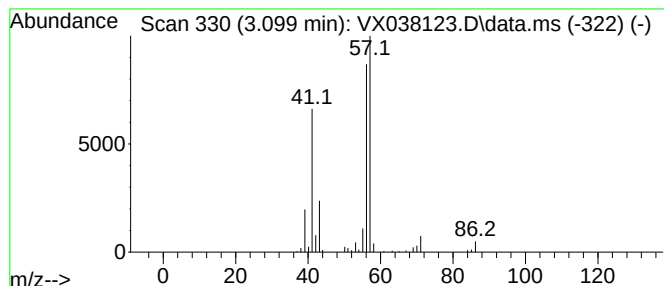
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Quant Title : SW846 8260

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

Peak Number 5 Pentane, 3-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.099	24.77 ug/l	139453	Pentafluorobenzene	5.555

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2			Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	64
3			Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	45
4			Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	43
5			Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100623\
Data File : VX038123.D
Acq On : 06 Oct 2023 12:14
Operator : JC/MD
Sample : 04723-02 20X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-5B

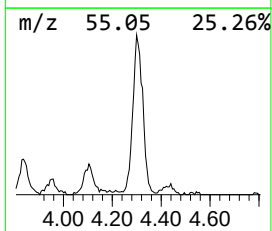
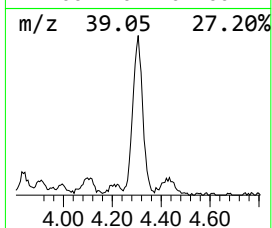
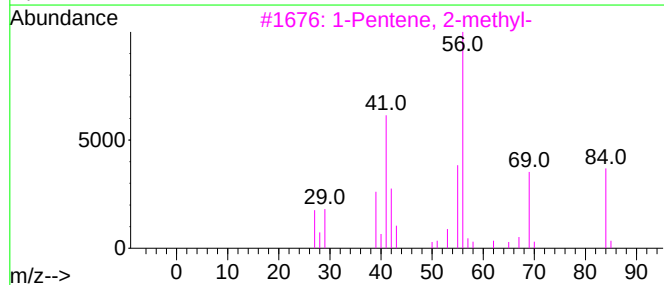
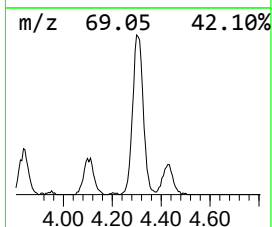
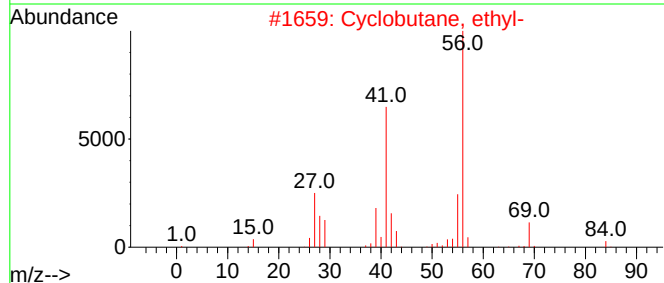
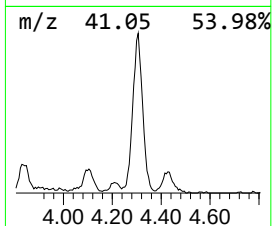
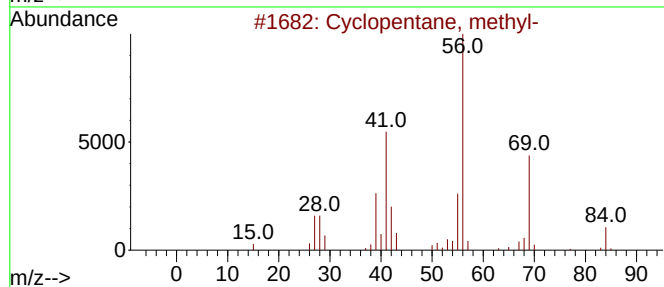
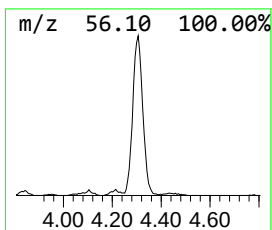
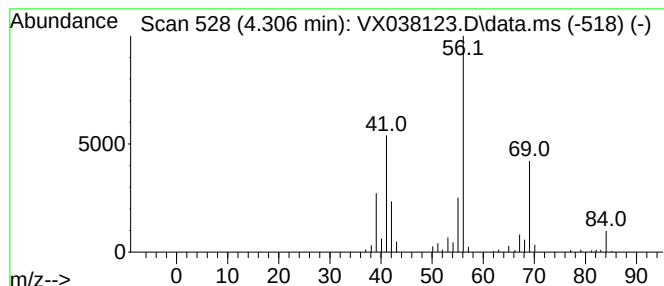
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Quant Title : SW846 8260

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

Peak Number 6 Cyclopentane, methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.306	34.60 ug/l	194777	Pentafluorobenzene	5.555

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	64
3			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	50
4			1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	50
5			Cyclobutane	56	C4H8	000287-23-0	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100623\
Data File : VX038123.D
Acq On : 06 Oct 2023 12:14
Operator : JC/MD
Sample : 04723-02 20X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-5B

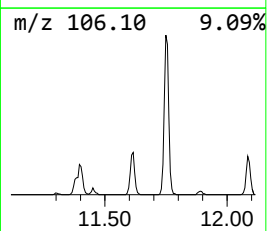
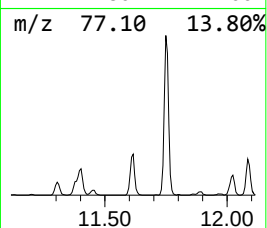
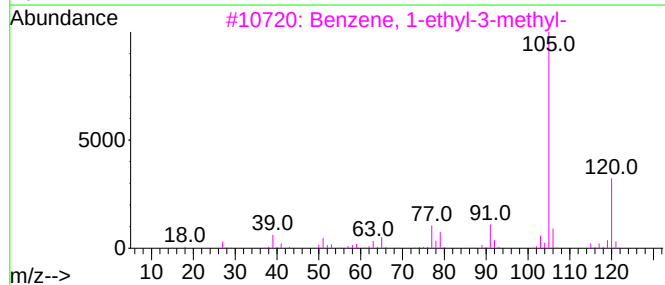
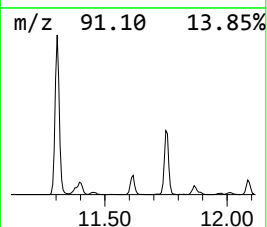
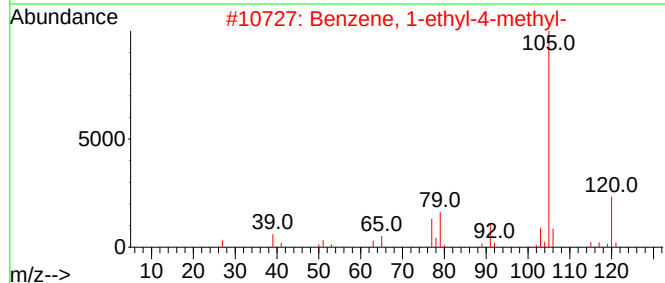
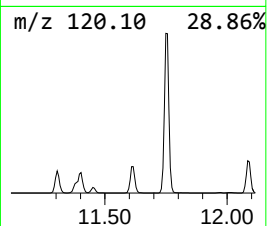
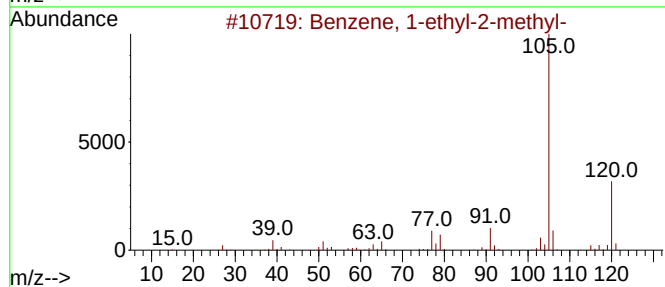
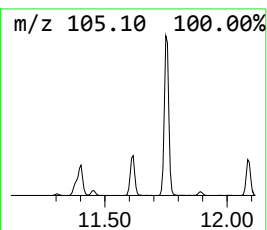
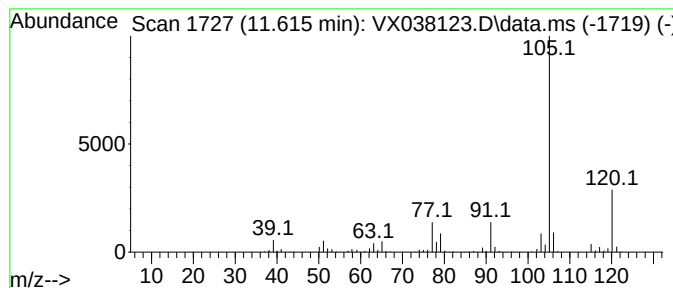
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.615	14.40 ug/l	193678	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	94
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100623\
Data File : VX038123.D
Acq On : 06 Oct 2023 12:14
Operator : JC/MD
Sample : 04723-02 20X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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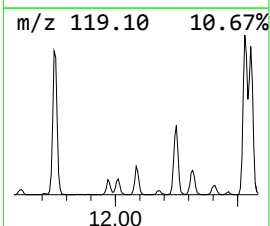
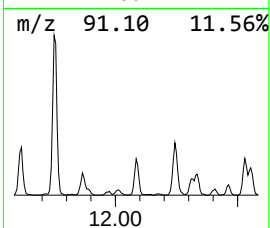
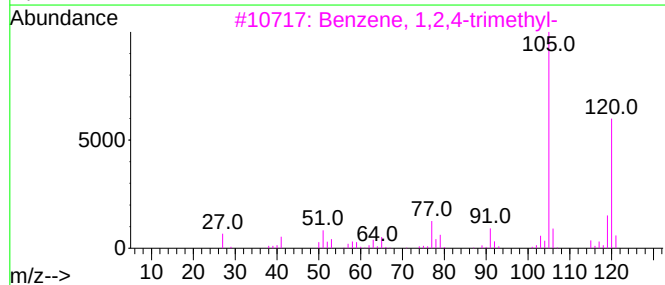
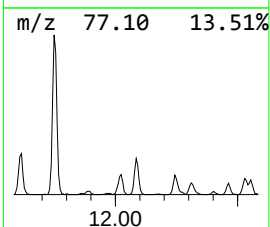
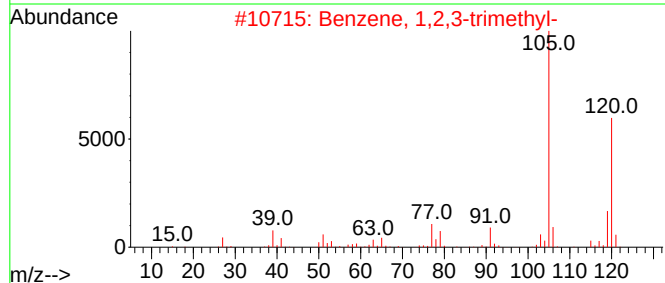
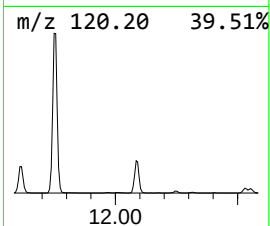
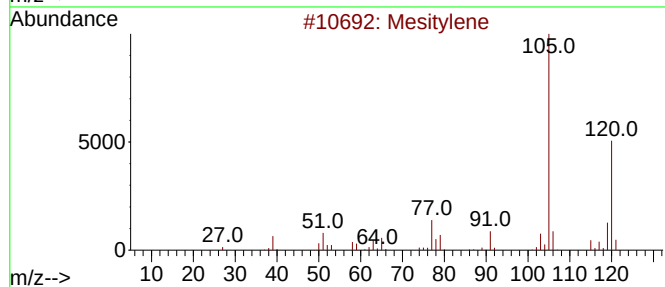
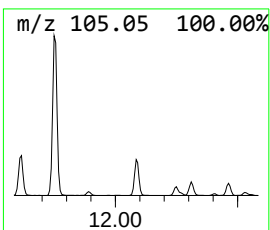
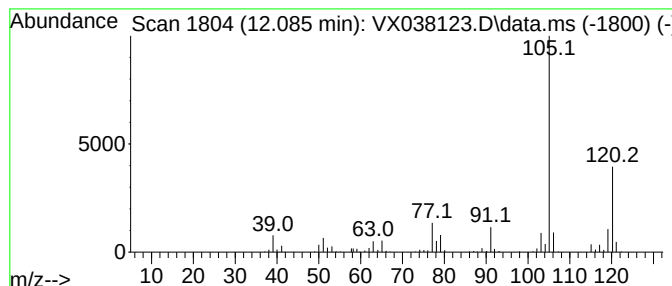
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.085	13.32 ug/l	179104	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	94
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100623\
Data File : VX038123.D
Acq On : 06 Oct 2023 12:14
Operator : JC/MD
Sample : 04723-02 20X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-5B

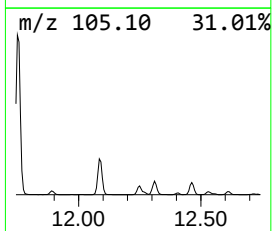
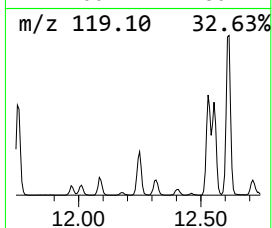
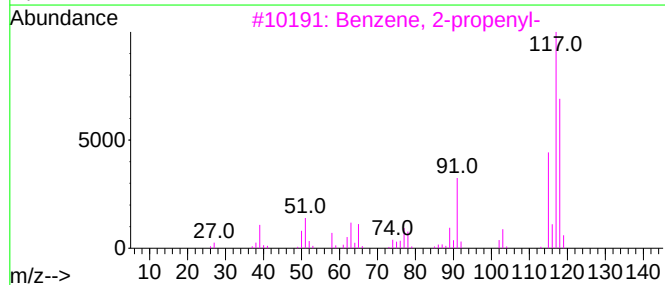
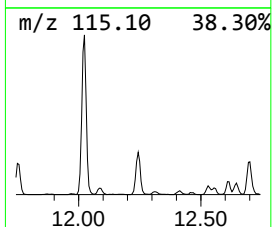
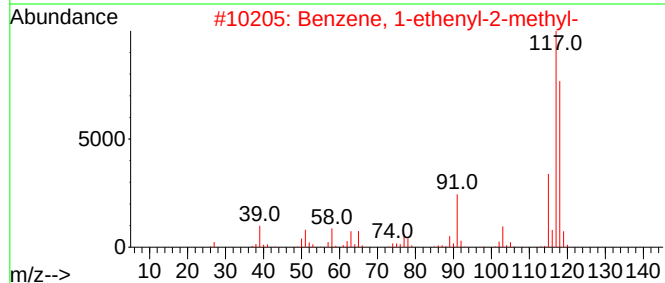
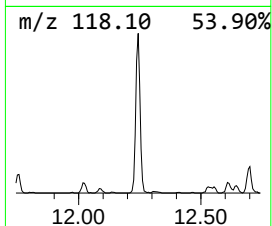
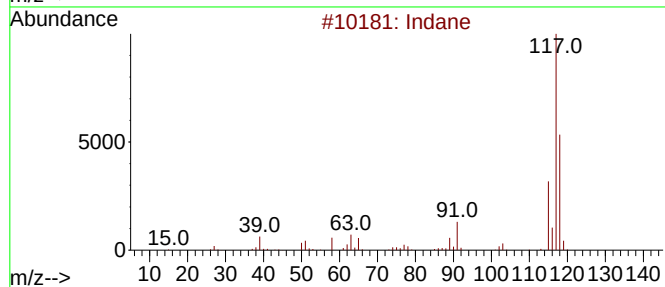
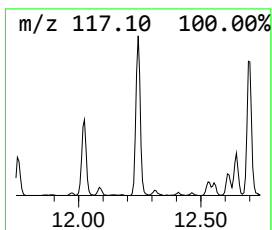
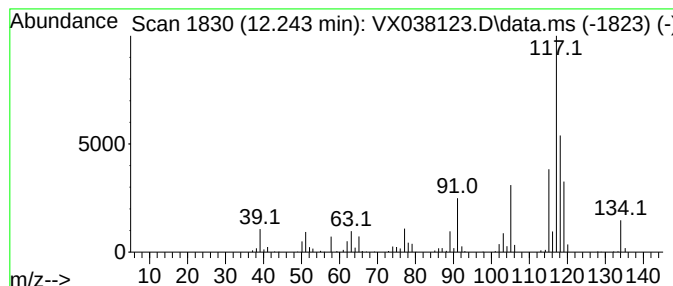
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Quant Title : SW846 8260

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

Peak Number 9 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.243	15.71 ug/l	211296	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	90
2			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	60
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	60
4			Benzene, 1-propenyl-	118	C9H10	000637-50-3	55
5			Benzene, cyclopropyl-	118	C9H10	000873-49-4	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100623\
Data File : VX038123.D
Acq On : 06 Oct 2023 12:14
Operator : JC/MD
Sample : 04723-02 20X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-5B

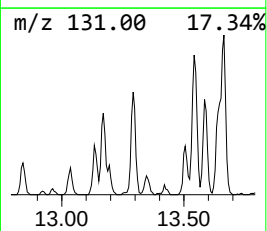
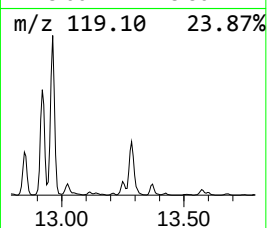
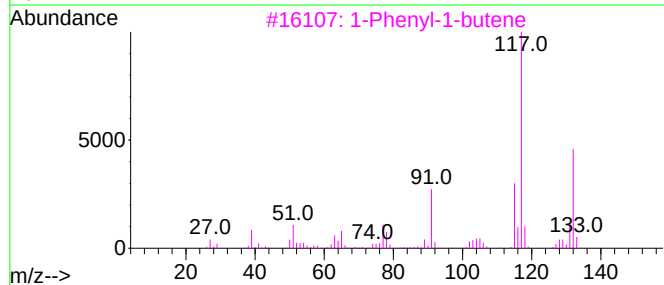
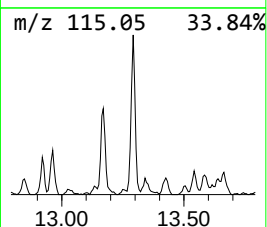
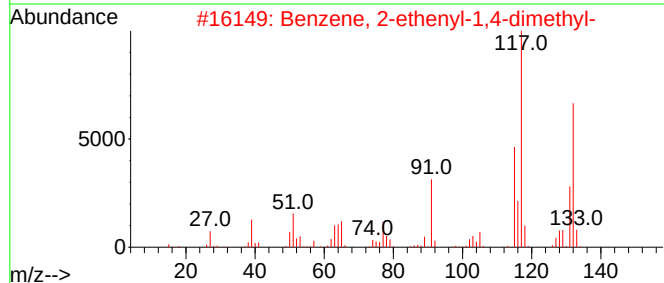
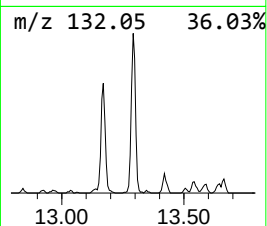
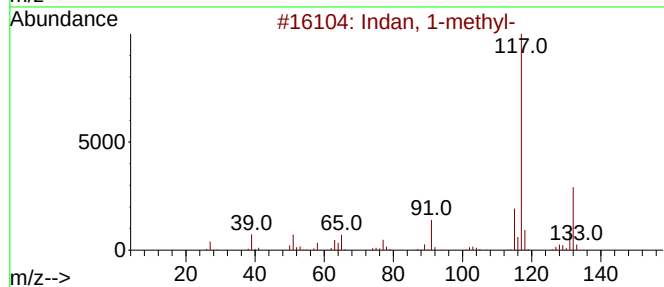
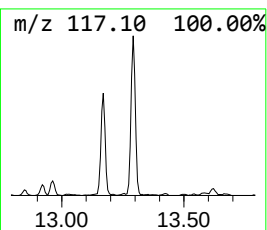
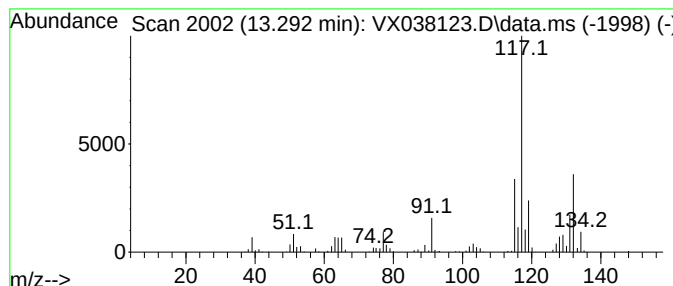
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Quant Title : SW846 8260

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

Peak Number 10 Indan, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	13.93 ug/l	187305	1,4-Dichlorobenzene-d4	12.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Indan, 1-methyl-	132	C10H12	000767-58-8	81
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	78
3		1-Phenyl-1-butene	132	C10H12	000824-90-8	74
4		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	74
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100623\
Data File : VX038123.D
Acq On : 06 Oct 2023 12:14
Operator : JC/MD
Sample : 04723-02 20X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-5B

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methyl-	1.745	61.2	ug/l	344421	1	5.555	281485	50.0
Pentane	1.952	17.4	ug/l	97880	1	5.555	281485	50.0
Butane, 2,3-dim...	2.782	15.1	ug/l	84978	1	5.555	281485	50.0
Pentane, 2-methyl-	2.824	17.6	ug/l	99194	1	5.555	281485	50.0
Pentane, 3-methyl-	3.099	24.8	ug/l	139453	1	5.555	281485	50.0
Cyclopentane, m...	4.306	34.6	ug/l	194777	1	5.555	281485	50.0
Benzene, 1-ethy...	11.615	14.4	ug/l	193678	4	12.024	672455	50.0
Benzene, 1,2,3-...	12.085	13.3	ug/l	179104	4	12.024	672455	50.0
Indane	12.243	15.7	ug/l	211296	4	12.024	672455	50.0
Indan, 1-methyl-	13.292	13.9	ug/l	187305	4	12.024	672455	50.0