

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120122\
 Data File : VX033155.D
 Acq On : 01 Dec 2022 12:05
 Operator : JC/MD
 Sample : N5815-13 10X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-10

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\82X112222W.M
 Title : SW846 8260

Signal : TIC: VX033155.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	43	47	52	rBV	298845	316849	5.28%	0.729%
2	1.746	103	108	125	rVB	1636773	2096871	34.94%	4.821%
3	1.953	136	142	155	rVB	397132	586174	9.77%	1.348%
4	2.215	180	185	194	rVB	78197	117989	1.97%	0.271%
5	2.782	270	278	281	rBV2	88177	195893	3.26%	0.450%
6	2.825	281	285	289	rVV	187216	368089	6.13%	0.846%
7	2.867	289	292	311	rVB4	88050	215934	3.60%	0.496%
8	3.105	320	331	347	rVB2	167138	398856	6.65%	0.917%
9	3.318	357	366	377	rVB2	27261	61562	1.03%	0.142%
10	3.446	379	387	399	rVB	47120	105522	1.76%	0.243%
11	3.727	427	433	443	rVB	52688	127551	2.13%	0.293%
12	3.837	443	451	457	rBV2	32774	79561	1.33%	0.183%
13	4.099	483	494	504	rVB2	35833	95068	1.58%	0.219%
14	4.215	504	513	518	rVV3	25219	70556	1.18%	0.162%
15	4.306	518	528	540	rVB	242944	699411	11.65%	1.608%
16	5.190	650	673	687	rVB3	36277	128660	2.14%	0.296%
17	5.385	693	705	712	rBV2	68910	202821	3.38%	0.466%
18	5.483	712	721	726	rVV4	58768	227812	3.80%	0.524%
19	5.550	726	732	739	rVV	149704	458881	7.65%	1.055%
20	5.617	739	743	762	rVB3	56412	173001	2.88%	0.398%
21	5.824	766	777	789	rBV2	48304	145735	2.43%	0.335%
22	5.958	789	799	806	rVV	71106	189884	3.16%	0.437%
23	6.037	806	812	822	rVV	57436	151732	2.53%	0.349%
24	6.159	823	832	839	rVV3	32096	111352	1.86%	0.256%
25	6.251	840	847	856	rVV3	62622	195150	3.25%	0.449%
26	6.361	856	865	876	rVB3	25993	79272	1.32%	0.182%
27	6.763	922	931	939	rBV	215560	544068	9.07%	1.251%
28	6.848	941	945	959	rVB5	33994	87834	1.46%	0.202%
29	7.379	1020	1032	1044	rBV5	69233	193307	3.22%	0.444%
30	7.982	1121	1131	1143	rVB	29894	72763	1.21%	0.167%
31	8.104	1143	1151	1154	rBV	50195	98873	1.65%	0.227%
32	8.507	1212	1217	1225	rVB2	32155	61246	1.02%	0.141%
33	8.647	1235	1240	1247	rVB	338256	530922	8.85%	1.221%
34	9.458	1365	1373	1377	rBV	50965	78180	1.30%	0.180%
35	10.055	1466	1471	1476	rBV	451129	598720	9.98%	1.377%
36	10.195	1488	1494	1500	rBV	1915394	2431955	40.52%	5.592%

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Integrator: RTE

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Stop Thrs : 0

Peak Location: TOP

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37	10.299	1505	1511	1518	rVB	2390653	3222595	53.70%	7.410%
38	10.640	1562	1567	1581	rVB	496615	636506	10.61%	1.464%
39	10.963	1614	1620	1626	rBV	277991	348864	5.81%	0.802%
40	11.079	1634	1639	1651	rBV	313949	408648	6.81%	0.940%
41	11.305	1667	1676	1683	rBV	820632	991874	16.53%	2.281%
42	11.378	1683	1688	1690	rBV	1918051	2475715	41.25%	5.692%
43	11.451	1696	1700	1712	rVB	1447849	1720539	28.67%	3.956%
44	11.610	1721	1726	1736	rBV	1548695	1952633	32.54%	4.490%
45	11.750	1744	1749	1755	rBV	4831911	6001262	100.00%	13.799%
46	11.866	1762	1768	1770	rBV	53800	72084	1.20%	0.166%
47	11.890	1770	1772	1777	rVB	65763	69692	1.16%	0.160%
48	11.969	1780	1785	1789	rBV	102074	122132	2.04%	0.281%
49	12.024	1789	1794	1799	rVB	528859	690785	11.51%	1.588%
50	12.085	1799	1804	1810	rBV	1650516	2026936	33.78%	4.660%
51	12.244	1824	1830	1838	rBV2	1335865	2059584	34.32%	4.736%
52	12.317	1838	1842	1849	rVB3	704482	1038464	17.30%	2.388%
53	12.408	1852	1857	1861	rBV2	65159	87459	1.46%	0.201%
54	12.463	1861	1866	1872	rVB	251784	307518	5.12%	0.707%
55	12.530	1872	1877	1879	rBV	403079	495772	8.26%	1.140%
56	12.555	1879	1881	1886	rVB	316232	333654	5.56%	0.767%
57	12.616	1886	1891	1894	rBV	802547	999198	16.65%	2.297%
58	12.646	1894	1896	1900	rVB	102954	105148	1.75%	0.242%
59	12.701	1900	1905	1913	rBV2	310629	486487	8.11%	1.119%
60	12.847	1924	1929	1934	rVB	192123	233879	3.90%	0.538%
61	12.920	1934	1941	1945	rBV	420855	525979	8.76%	1.209%
62	12.963	1945	1948	1954	rVB	468639	536810	8.94%	1.234%
63	13.024	1954	1958	1960	rBV	47979	65803	1.10%	0.151%
64	13.170	1977	1982	1986	rVB	350648	422017	7.03%	0.970%
65	13.256	1992	1996	1998	rBV2	48810	72823	1.21%	0.167%
66	13.292	1998	2002	2007	rVB2	696861	911386	15.19%	2.096%
67	13.372	2012	2015	2019	rVV	51711	65638	1.09%	0.151%
68	13.426	2019	2024	2031	rVB	85893	115784	1.93%	0.266%
69	13.579	2046	2049	2056	rVV4	89842	159416	2.66%	0.367%
70	13.664	2056	2063	2069	rVB2	56712	130824	2.18%	0.301%
71	13.774	2075	2081	2088	rBV	516346	639258	10.65%	1.470%
72	14.073	2126	2130	2135	rBV	54063	68528	1.14%	0.158%
73	14.213	2149	2153	2159	rBV	45205	76075	1.27%	0.175%
74	14.640	2213	2223	2234	rBV	226739	303399	5.06%	0.698%
75	14.780	2241	2246	2255	rVB	172865	212719	3.54%	0.489%

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Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Title : SW846 8260

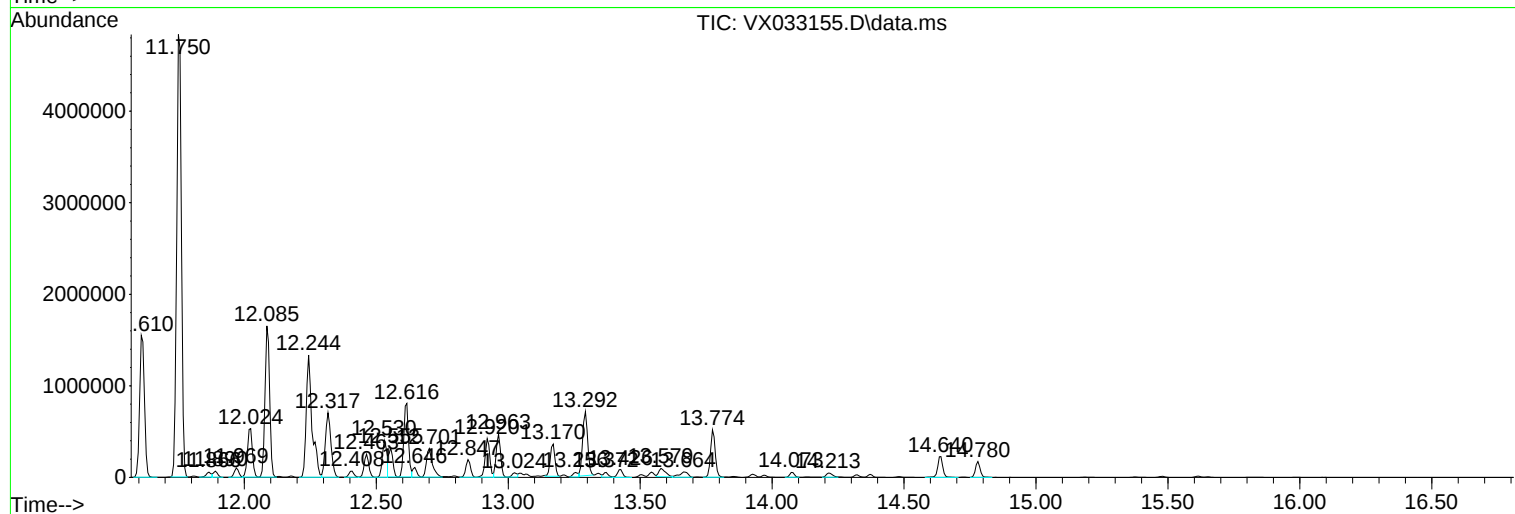
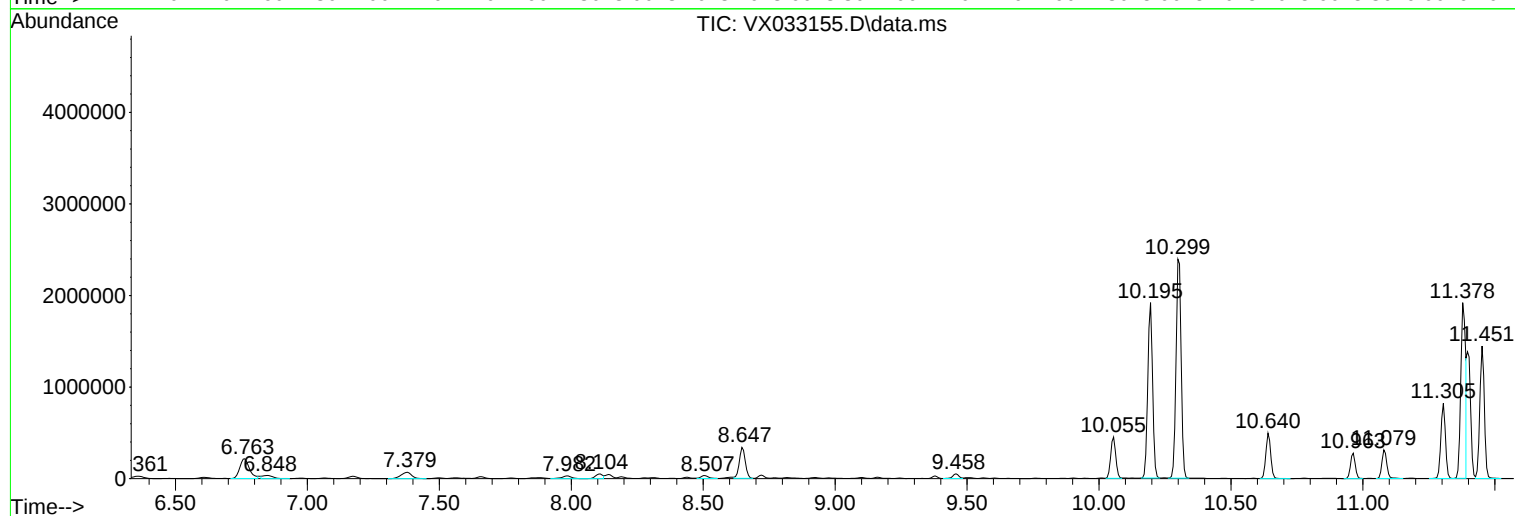
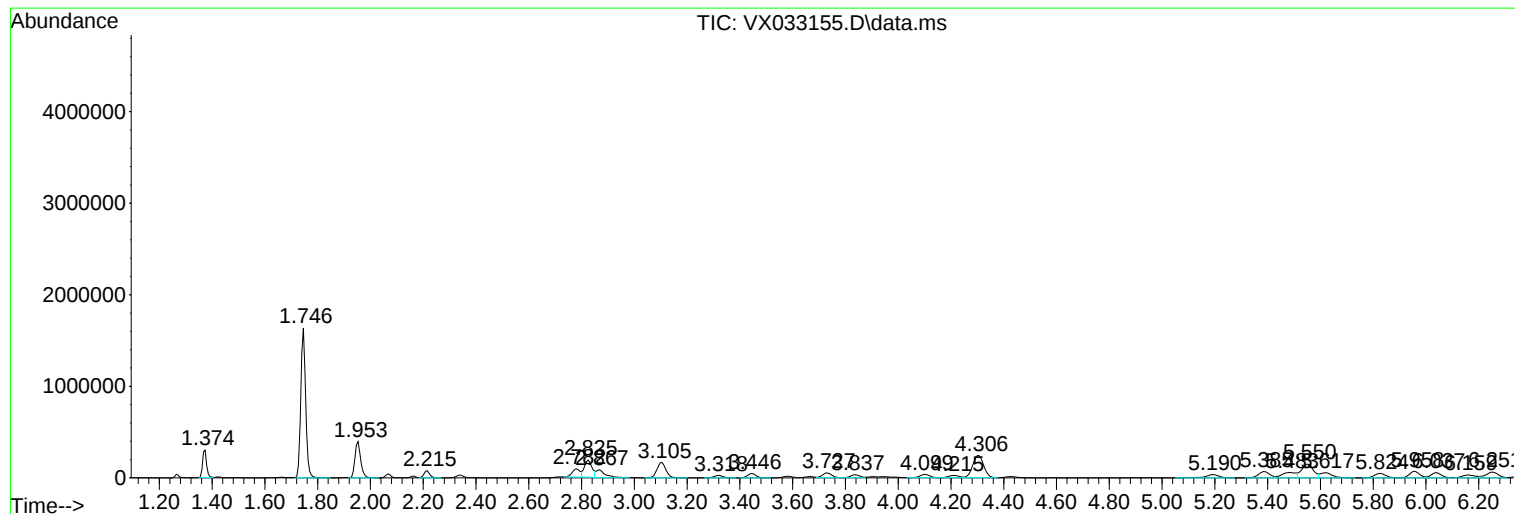
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Instrument :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

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



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Instrument :
MSVOA_X
ClientSampleId :
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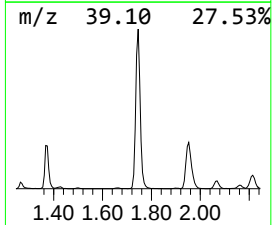
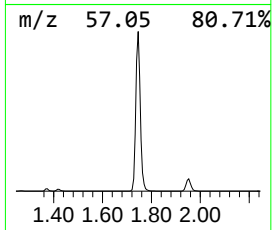
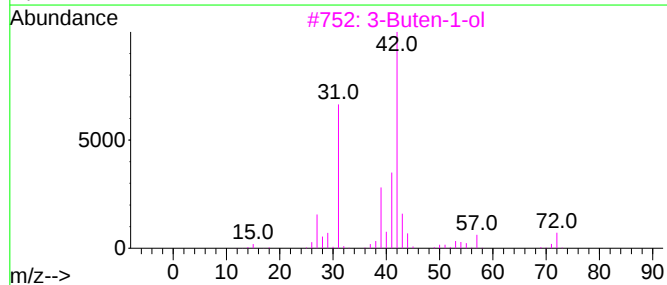
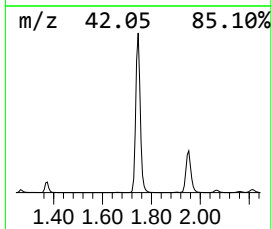
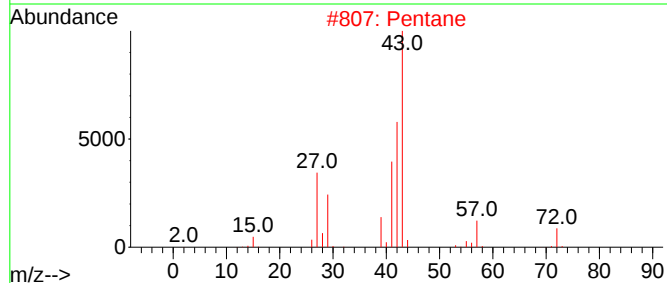
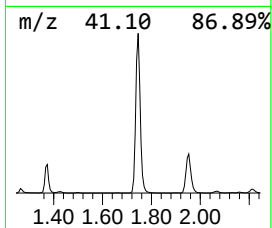
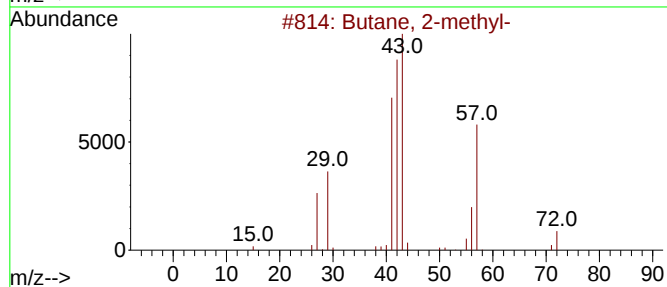
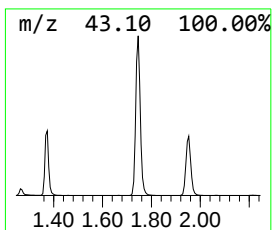
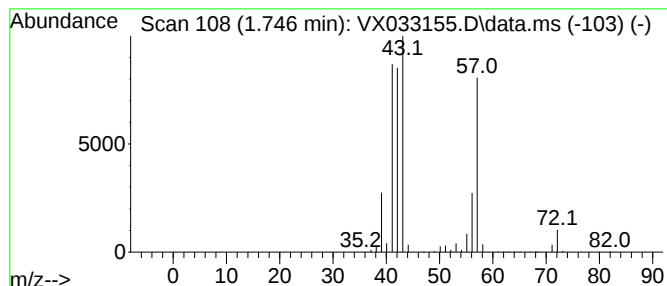
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

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.
1.746	228.48 ug/l	2096870	Pentafluorobenzene	5.550

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methyl-	72	C5H12	000078-78-4	87
2			Pentane	72	C5H12	000109-66-0	42
3			3-Buten-1-ol	72	C4H8O	000627-27-0	25
4			Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	10
5			Butane, 1-chloro-2-methyl-	106	C5H11Cl	000616-13-7	10



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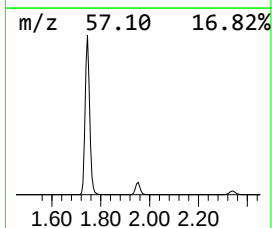
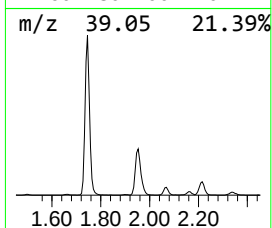
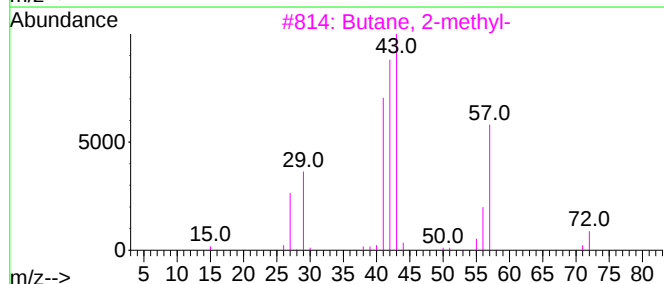
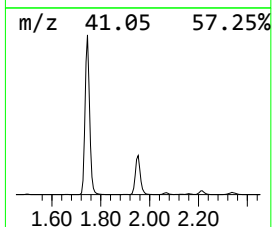
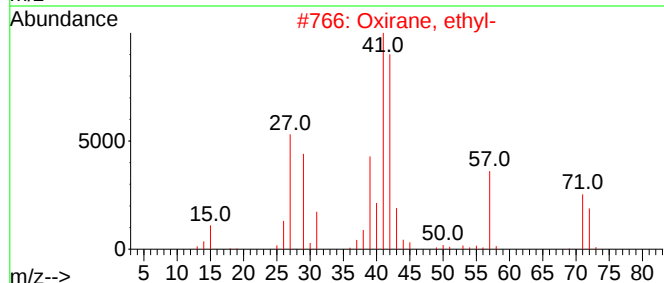
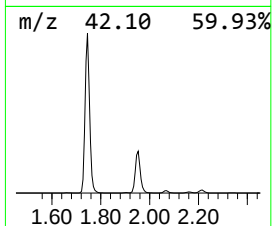
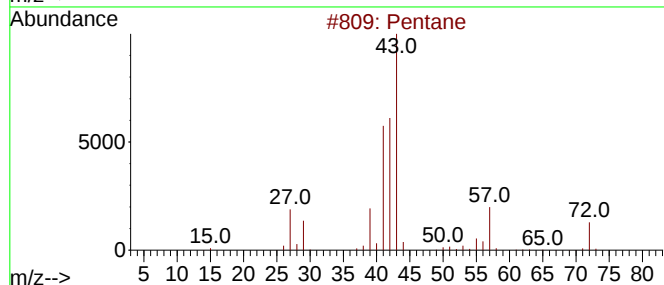
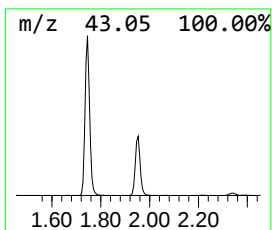
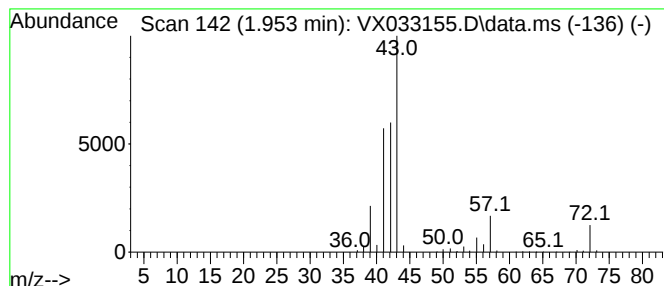
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Pentane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.953	63.87 ug/l	586174	Pentafluorobenzene	5.550

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			Butane, 2-methyl-	72	C5H12	000078-78-4	50
4			2-Propanone, hydrazone	72	C3H8N2	005281-20-9	10
5			1-Propene, 2-methoxy-	72	C4H8O	000116-11-0	9



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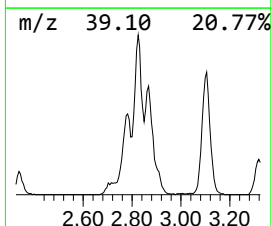
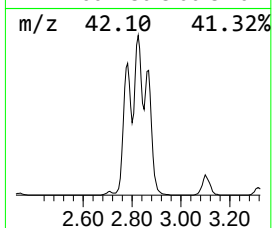
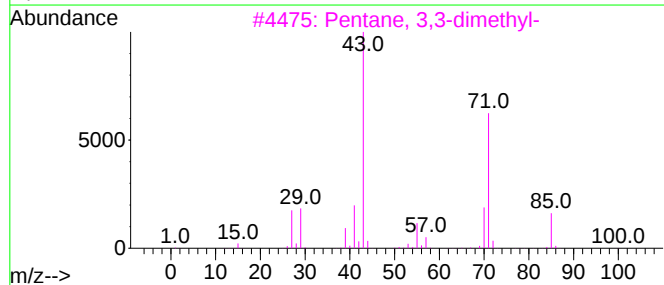
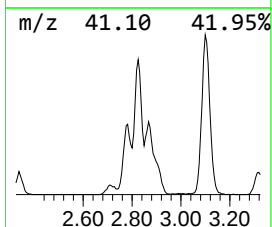
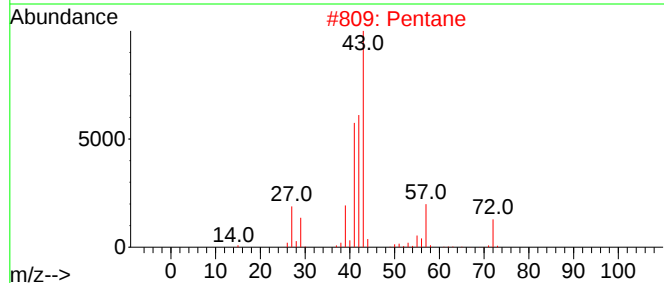
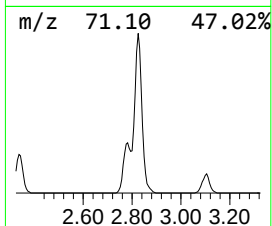
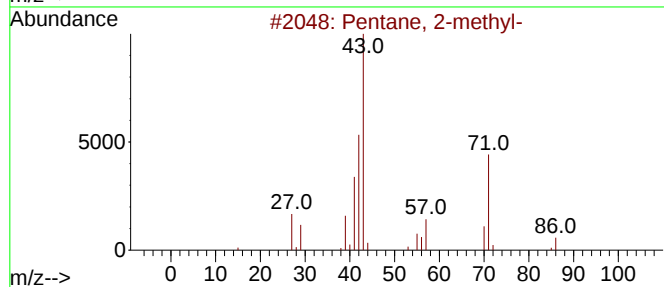
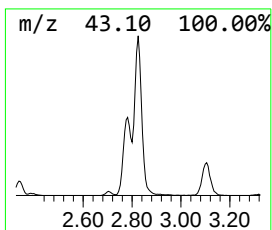
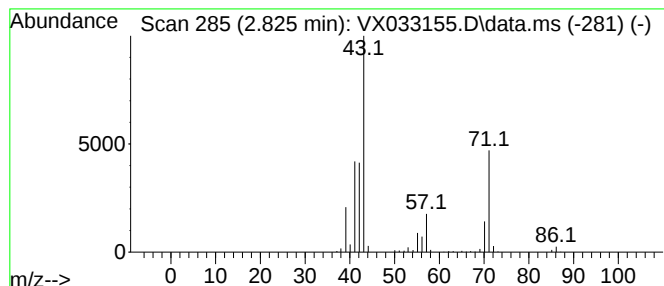
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Quant Title : SW846 8260

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

Peak Number 3 Pentane, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.825	40.11 ug/l	368089	Pentafluorobenzene	5.550

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2-methyl-	86	C6H14	000107-83-5	91
2			Pentane	72	C5H12	000109-66-0	46
3			Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	38
4			Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	36
5			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	28



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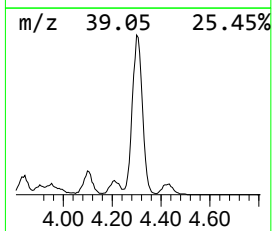
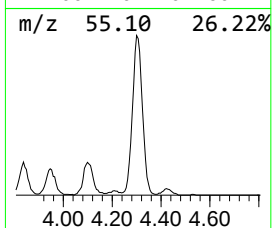
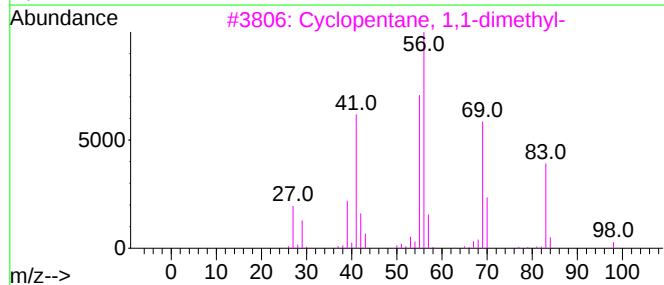
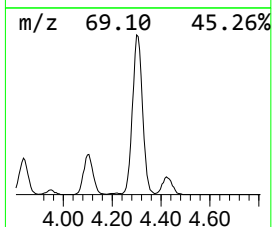
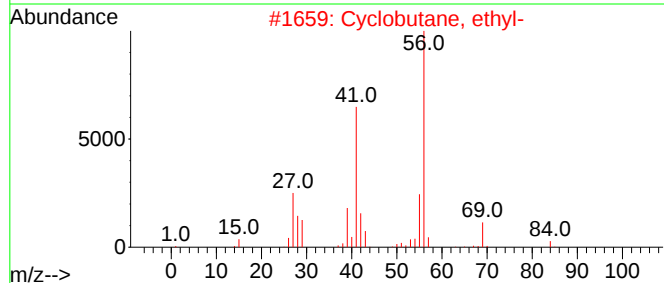
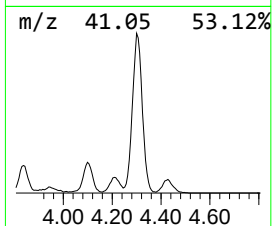
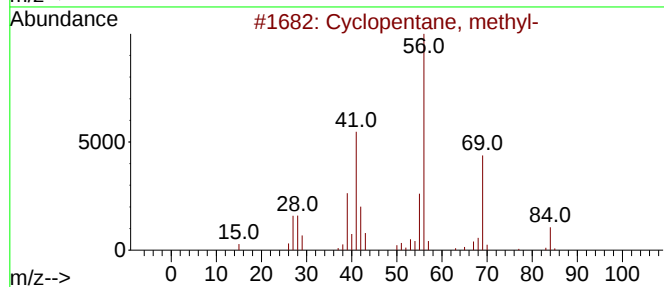
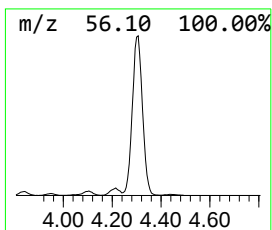
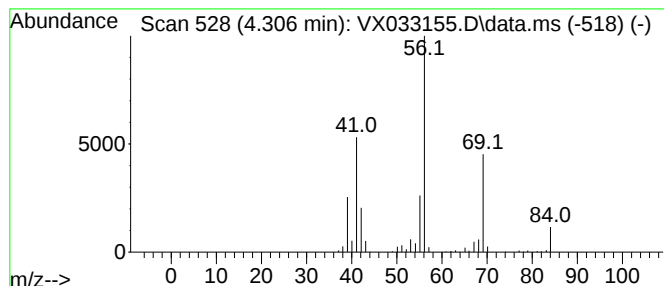
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Quant Title : SW846 8260

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

Peak Number 4 Cyclopentane, methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.306	76.21 ug/l	699411	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	72
3			Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	72
4			1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64
5			1-Hexene	84	C6H12	000592-41-6	52



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120122\
Data File : VX033155.D
Acq On : 01 Dec 2022 12:05
Operator : JC/MD
Sample : N5815-13 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10

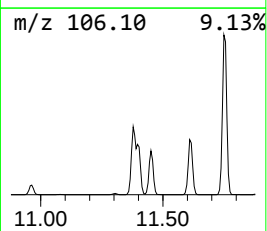
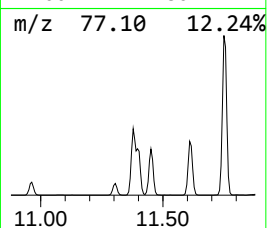
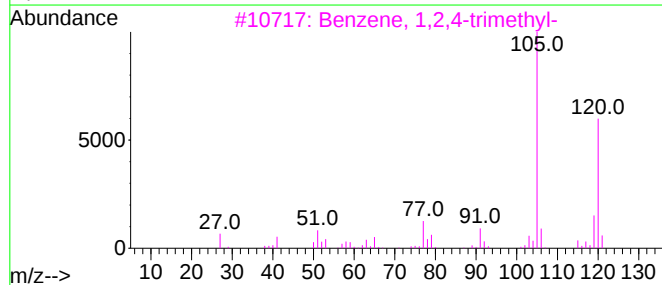
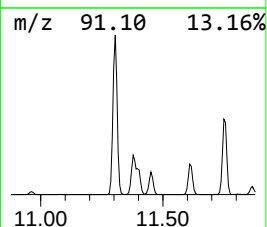
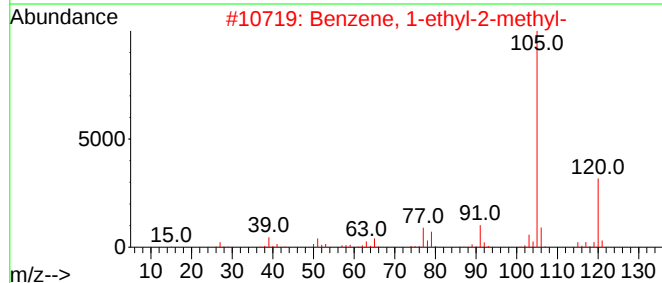
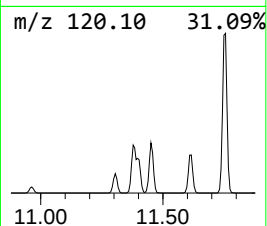
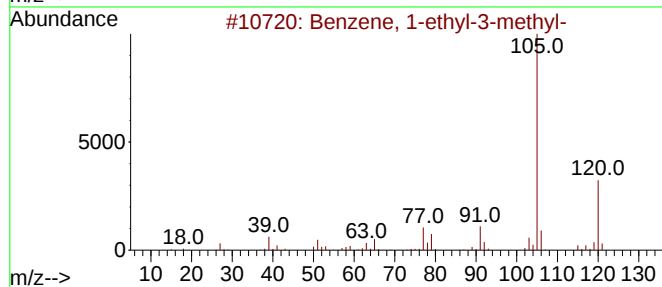
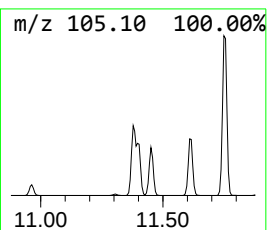
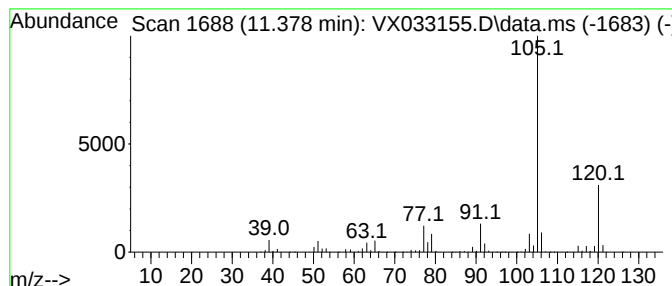
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-ethyl-3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.378	179.20 ug/l	2475720	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	95
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			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120122\
Data File : VX033155.D
Acq On : 01 Dec 2022 12:05
Operator : JC/MD
Sample : N5815-13 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10

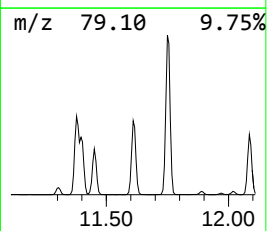
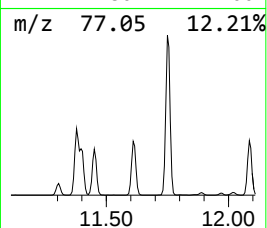
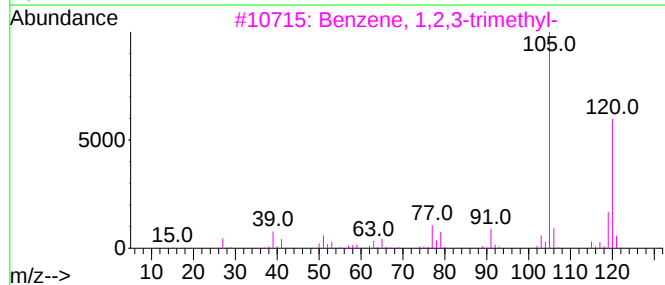
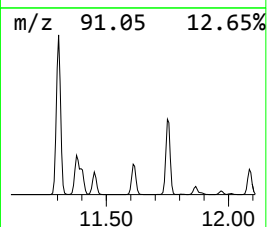
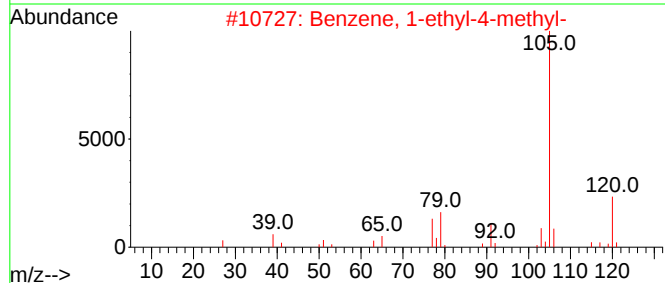
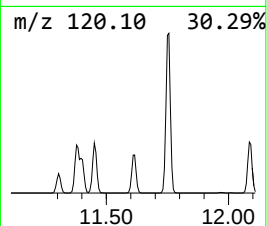
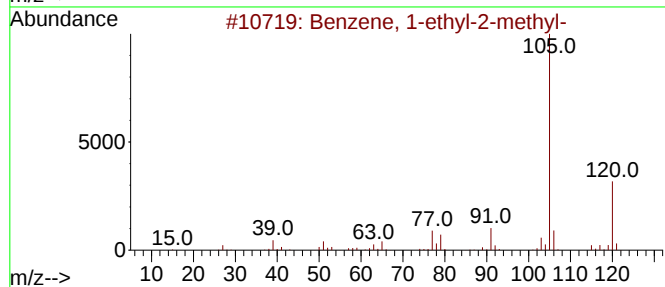
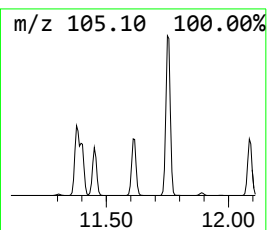
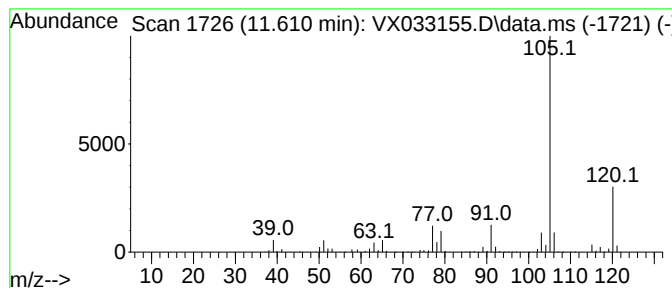
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethyl-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.610	141.33 ug/l	1952630	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	93
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	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\VX120122\
Data File : VX033155.D
Acq On : 01 Dec 2022 12:05
Operator : JC/MD
Sample : N5815-13 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10

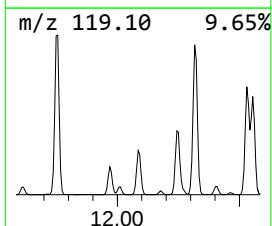
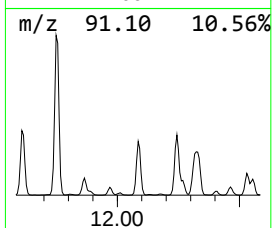
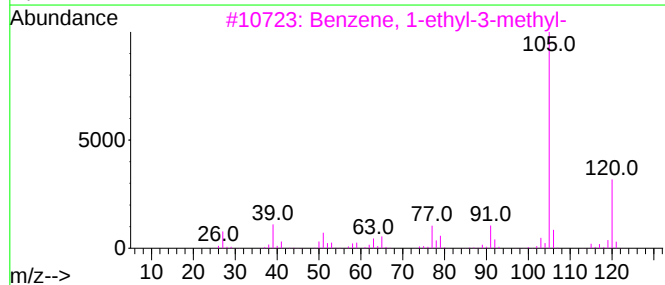
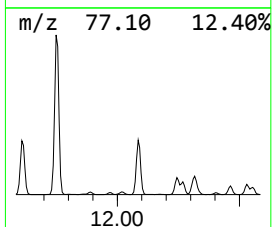
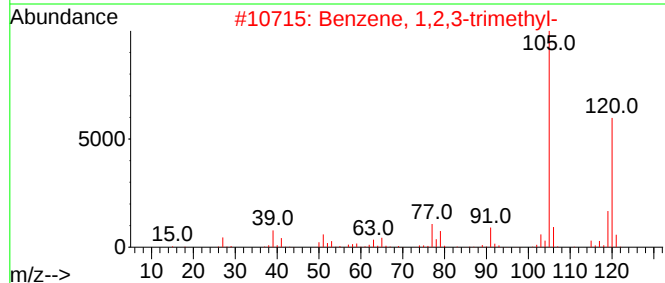
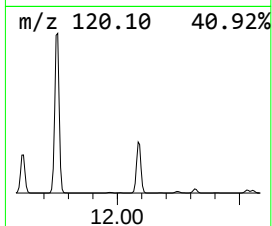
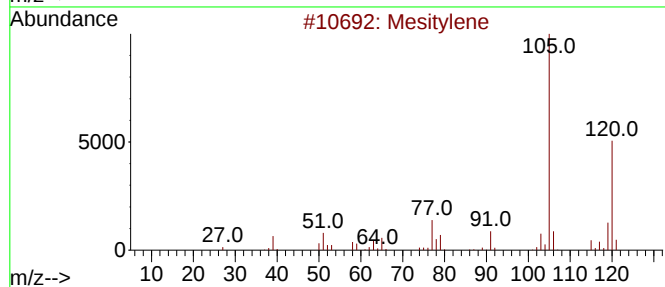
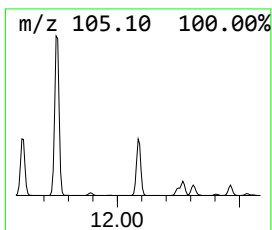
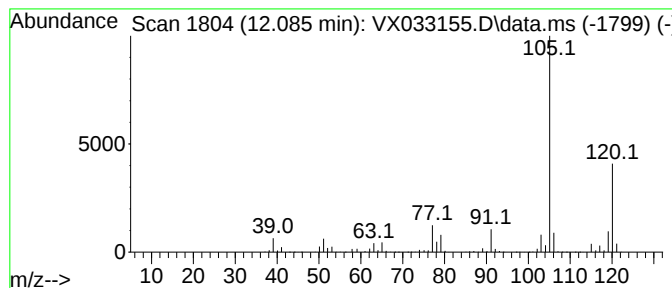
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,3-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.085	146.71 ug/l	2026940	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-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120122\
Data File : VX033155.D
Acq On : 01 Dec 2022 12:05
Operator : JC/MD
Sample : N5815-13 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10

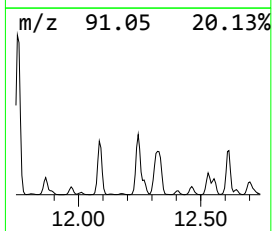
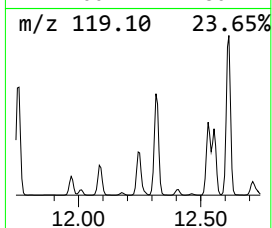
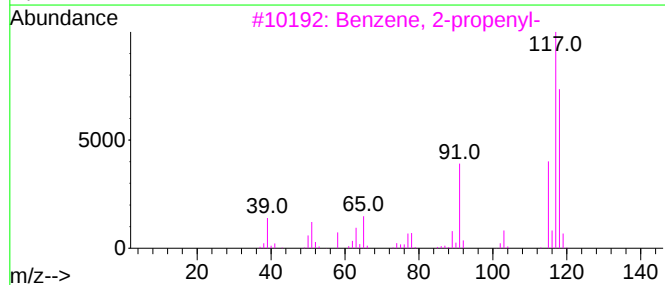
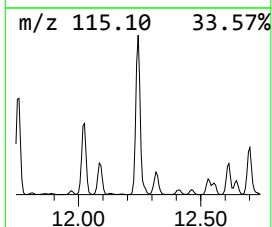
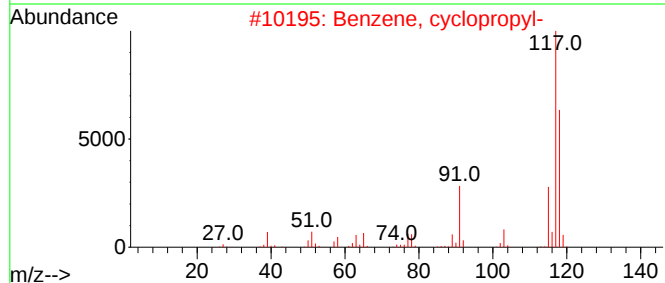
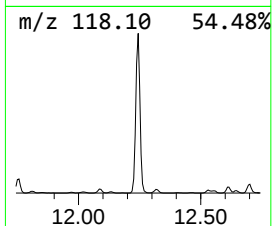
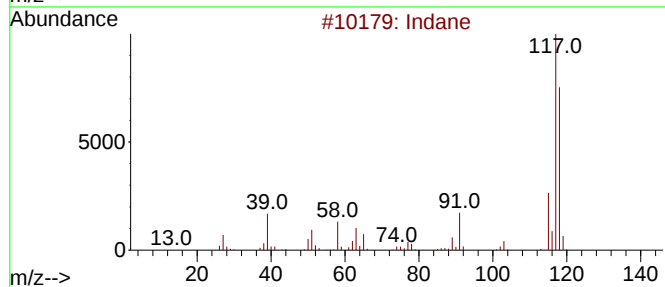
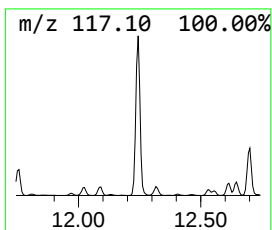
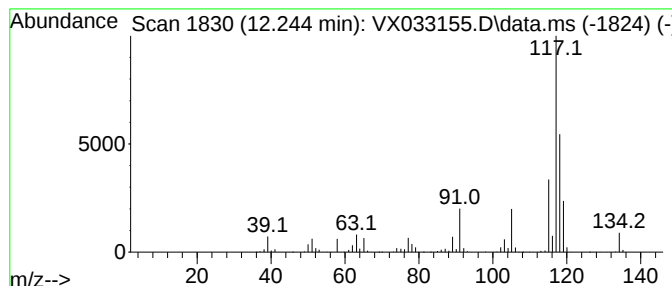
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Quant Title : SW846 8260

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

Peak Number 8 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	149.08 ug/l	2059580	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	70
2			Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	70
4			Benzene, 1-propenyl-	118	C9H10	000637-50-3	70
5			Deltacyclene	118	C9H10	007785-10-6	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120122\
Data File : VX033155.D
Acq On : 01 Dec 2022 12:05
Operator : JC/MD
Sample : N5815-13 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

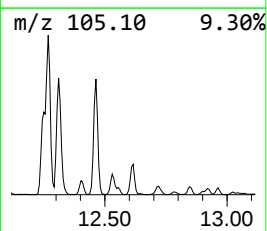
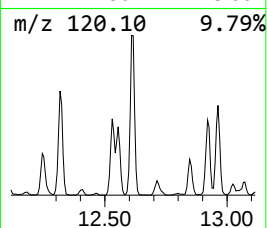
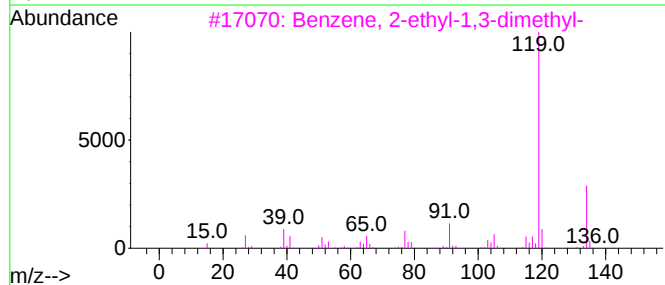
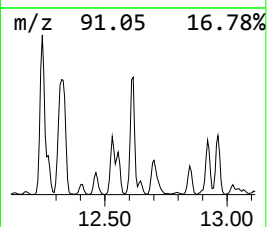
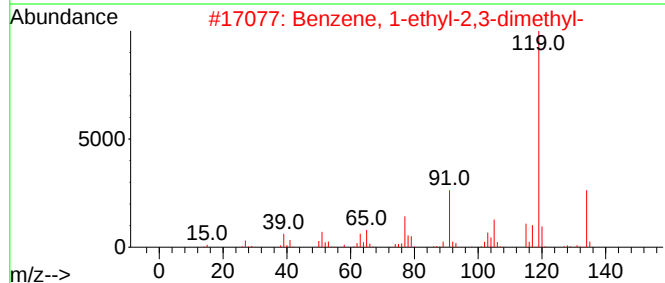
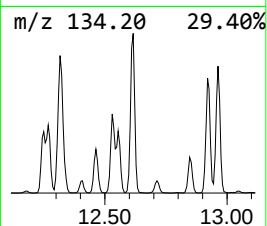
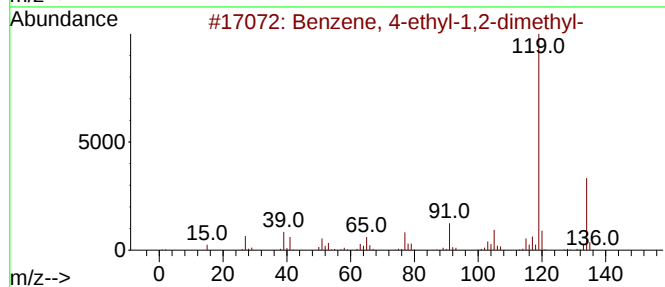
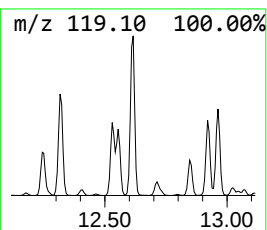
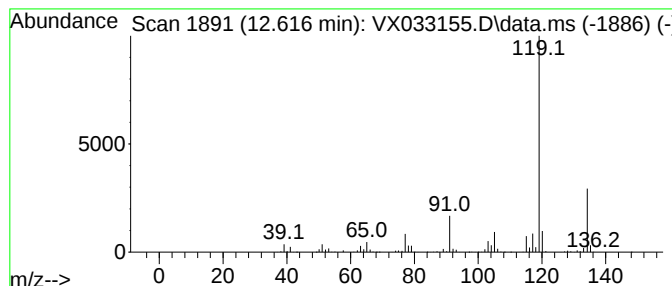
Instrument :
MSVOA_X
ClientSampleId :
MW-10

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD			R.T.
12.616	72.32 ug/l	999198	1,4-Dichlorobenzene-d4			12.024
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-		134	C10H14	000934-80-5	95
2	Benzene, 1-ethyl-2,3-dimethyl-		134	C10H14	000933-98-2	95
3	Benzene, 2-ethyl-1,3-dimethyl-		134	C10H14	002870-04-4	95
4	o-Cymene		134	C10H14	000527-84-4	95
5	Benzene, 1-ethyl-2,4-dimethyl-		134	C10H14	000874-41-9	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120122\
Data File : VX033155.D
Acq On : 01 Dec 2022 12:05
Operator : JC/MD
Sample : N5815-13 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10

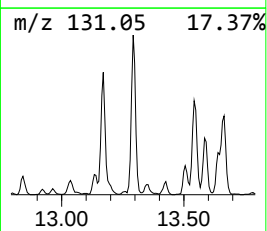
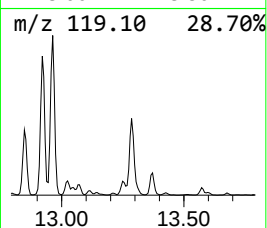
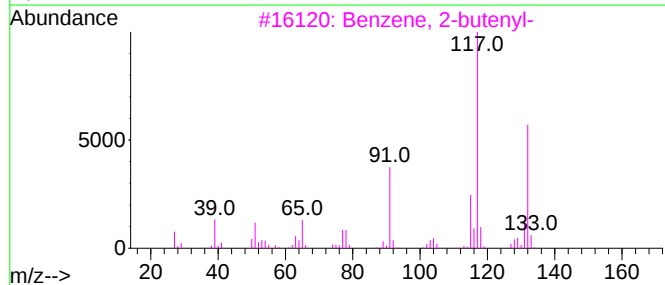
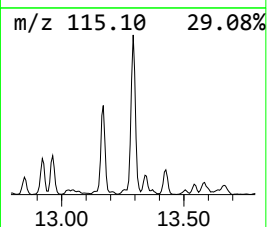
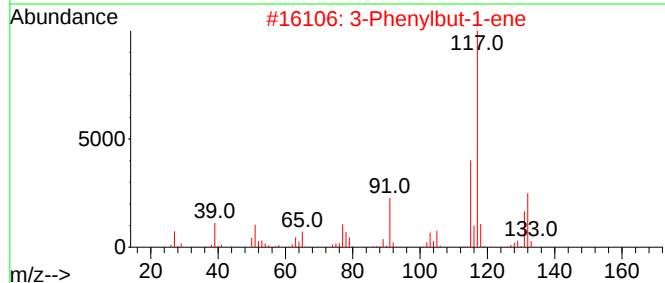
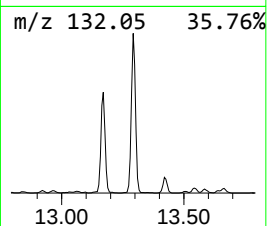
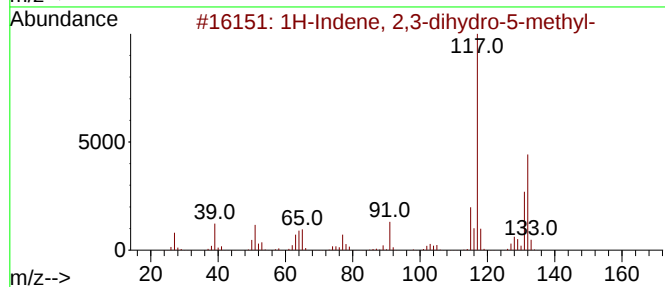
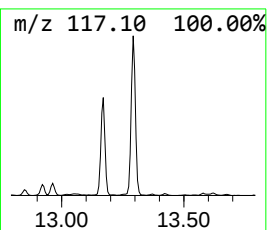
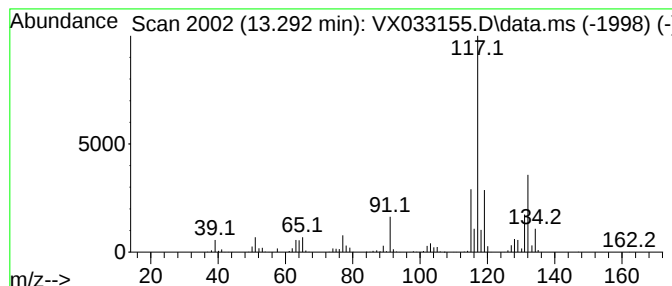
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Quant Title : SW846 8260

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

Peak Number 10 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	65.97 ug/l	911386	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		3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
3		Benzene, 2-butenyl-	132	C10H12	001560-06-1	81
4		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	76
5		Indan, 1-methyl-	132	C10H12	000767-58-8	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120122\
Data File : VX033155.D
Acq On : 01 Dec 2022 12:05
Operator : JC/MD
Sample : N5815-13 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.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-	1.746	228.5	ug/l	2096870	1	5.550	458881	50.0
Pentane	1.953	63.9	ug/l	586174	1	5.550	458881	50.0
Pentane, 2-methyl-	2.825	40.1	ug/l	368089	1	5.550	458881	50.0
Cyclopentane, m...	4.306	76.2	ug/l	699411	1	5.550	458881	50.0
Benzene, 1-ethy...	11.378	179.2	ug/l	2475720	4	12.024	690785	50.0
Benzene, 1-ethy...	11.610	141.3	ug/l	1952630	4	12.024	690785	50.0
Benzene, 1,2,3-...	12.085	146.7	ug/l	2026940	4	12.024	690785	50.0
Indane	12.244	149.1	ug/l	2059580	4	12.024	690785	50.0
Benzene, 4-ethy...	12.616	72.3	ug/l	999198	4	12.024	690785	50.0
1H-Indene, 2,3-...	13.292	66.0	ug/l	911386	4	12.024	690785	50.0