

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120122\
 Data File : VX033156.D
 Acq On : 01 Dec 2022 12:28
 Operator : JC/MD
 Sample : N5815-25
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-15

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: VX033156.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.264	26	29	32	rBV	238076	234944	1.02%	0.111%
2	1.374	43	47	54	rBV	2110245	2379351	10.38%	1.123%
3	1.752	103	109	127	rVB	8994570	11828896	51.60%	5.581%
4	1.959	138	143	156	rVB	3069670	4414889	19.26%	2.083%
5	2.069	156	161	168	rVV	447181	642235	2.80%	0.303%
6	2.215	181	185	193	rVB	188740	296625	1.29%	0.140%
7	2.343	200	206	210	rBV	245468	471342	2.06%	0.222%
8	2.386	210	213	229	rVB	203653	357828	1.56%	0.169%
9	2.782	271	278	281	rVV	860564	1808153	7.89%	0.853%
10	2.831	281	286	290	rVV	2099861	4538480	19.80%	2.141%
11	2.867	290	292	308	rVB2	1034338	1994888	8.70%	0.941%
12	3.105	320	331	352	rVB	1551235	3442197	15.02%	1.624%
13	3.318	355	366	377	rBV2	144456	336755	1.47%	0.159%
14	3.447	377	387	400	rBV	748132	1660202	7.24%	0.783%
15	3.587	400	410	416	rBV2	120262	305086	1.33%	0.144%
16	3.672	416	424	429	rVV	125393	302414	1.32%	0.143%
17	3.733	429	434	443	rVB	168556	388675	1.70%	0.183%
18	3.837	443	451	457	rBV	91441	233082	1.02%	0.110%
19	3.946	466	469	482	rVB2	82606	230166	1.00%	0.109%
20	4.105	482	495	504	rVB2	117809	324038	1.41%	0.153%
21	4.215	504	513	519	rBV	207312	612147	2.67%	0.289%
22	4.306	519	528	543	rVB	2931695	8372203	36.52%	3.950%
23	4.556	560	569	584	rVB	122899	352732	1.54%	0.166%
24	5.385	690	705	711	rBV4	93349	317951	1.39%	0.150%
25	5.483	711	721	736	rVV5	601102	2935057	12.80%	1.385%
26	5.623	737	744	765	rVB	430420	1472358	6.42%	0.695%
27	5.830	765	778	791	rBV2	432486	1296535	5.66%	0.612%
28	6.038	804	812	822	rVV	422395	1140134	4.97%	0.538%
29	6.159	822	832	839	rVV2	293378	987877	4.31%	0.466%
30	6.257	840	848	857	rVV2	630687	1970155	8.59%	0.930%
31	6.361	857	865	876	rVB	240809	691558	3.02%	0.326%
32	6.617	893	907	916	rBV4	283143	871620	3.80%	0.411%
33	6.763	923	931	937	rBV	232001	588889	2.57%	0.278%
34	6.848	938	945	956	rVB4	150888	495234	2.16%	0.234%
35	7.379	1018	1032	1040	rBV4	830314	2272420	9.91%	1.072%
36	7.659	1071	1078	1089	rVB2	252142	561368	2.45%	0.265%

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37	7.891	1105	1116	1123	rBV4	103830	321617	1.40%	0.152%
38	7.982	1123	1131	1143	rVB	294332	648912	2.83%	0.306%
39	8.104	1143	1151	1159	rBV	425554	894450	3.90%	0.422%
40	8.190	1159	1165	1174	rVB2	122297	242043	1.06%	0.114%
41	8.433	1200	1205	1213	rBV2	192676	364460	1.59%	0.172%
42	8.507	1213	1217	1224	rVB3	145279	251916	1.10%	0.119%
43	8.610	1224	1234	1236	rBV5	153428	321740	1.40%	0.152%
44	8.647	1236	1240	1246	rVB	378291	615653	2.69%	0.290%
45	8.714	1246	1251	1257	rBV3	304717	530610	2.31%	0.250%
46	8.805	1257	1266	1270	rBV	202843	437221	1.91%	0.206%
47	10.055	1466	1471	1475	rVB	462226	639655	2.79%	0.302%
48	10.201	1488	1495	1501	rBV	16764830	22922206	100.00%	10.815%
49	10.311	1503	1513	1518	rVB	14212539	19028597	83.01%	8.978%
50	10.640	1562	1567	1578	rVB	505930	704454	3.07%	0.332%
51	10.963	1613	1620	1626	rBV	1989885	2489287	10.86%	1.174%
52	11.079	1634	1639	1643	rBV	347362	466273	2.03%	0.220%
53	11.305	1668	1676	1682	rBV	6070667	7403599	32.30%	3.493%
54	11.402	1682	1692	1696	rVV	7791984	9526697	41.56%	4.495%
55	11.451	1696	1700	1712	rVB	2024660	2473751	10.79%	1.167%
56	11.616	1721	1727	1734	rVB	2783138	3579857	15.62%	1.689%
57	11.756	1744	1750	1755	rVB	16173038	19561630	85.34%	9.229%
58	11.866	1761	1768	1770	rBV	347029	456828	1.99%	0.216%
59	11.890	1770	1772	1777	rVB	450047	476041	2.08%	0.225%
60	11.969	1781	1785	1789	rVV	251500	307826	1.34%	0.145%
61	12.018	1789	1793	1799	rVB2	615379	903858	3.94%	0.426%
62	12.091	1799	1805	1810	rBV	4577164	5724630	24.97%	2.701%
63	12.244	1824	1830	1837	rBV	5987583	7624607	33.26%	3.597%
64	12.317	1837	1842	1850	rVB3	2475980	4186041	18.26%	1.975%
65	12.408	1852	1857	1861	rBV	348852	446324	1.95%	0.211%
66	12.463	1861	1866	1872	rVB	1421754	1671060	7.29%	0.788%
67	12.530	1872	1877	1880	rBV	2052752	2870879	12.52%	1.354%
68	12.555	1880	1881	1886	rVB	1432666	1140485	4.98%	0.538%
69	12.616	1886	1891	1894	rBV	3820932	4441842	19.38%	2.096%
70	12.646	1894	1896	1900	rVB	510448	529126	2.31%	0.250%
71	12.701	1900	1905	1913	rBV2	1469055	2253674	9.83%	1.063%
72	12.799	1913	1921	1925	rBV	293460	391004	1.71%	0.184%
73	12.847	1925	1929	1934	rVB	931309	1128449	4.92%	0.532%
74	12.920	1934	1941	1944	rBV	1821617	2324663	10.14%	1.097%
75	12.963	1944	1948	1954	rVB	1902765	2428993	10.60%	1.146%
76	13.024	1954	1958	1960	rBV	249836	317947	1.39%	0.150%

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77	13.170	1978	1982	1986	rVB	1214240	1430309	6.24%	0.675%
78	13.256	1992	1996	1998	rBV2	256859	364501	1.59%	0.172%
79	13.292	1998	2002	2008	rVV2	2618249	3547615	15.48%	1.674%
80	13.372	2012	2015	2019	rVB	225795	262805	1.15%	0.124%
81	13.420	2019	2023	2032	rVB	285891	393530	1.72%	0.186%
82	13.542	2040	2043	2046	rVV	208066	260106	1.13%	0.123%
83	13.579	2046	2049	2056	rVV4	426258	885657	3.86%	0.418%
84	13.664	2060	2063	2068	rVB2	226798	393588	1.72%	0.186%
85	13.774	2076	2081	2086	rBV	3662709	4643850	20.26%	2.191%
86	13.823	2086	2089	2094	rVB2	733769	902670	3.94%	0.426%
87	14.213	2149	2153	2159	rBV2	173513	272317	1.19%	0.128%
88	14.640	2218	2223	2228	rVB	2262052	2855015	12.46%	1.347%
89	14.780	2241	2246	2255	rVB	1239660	1565432	6.83%	0.739%

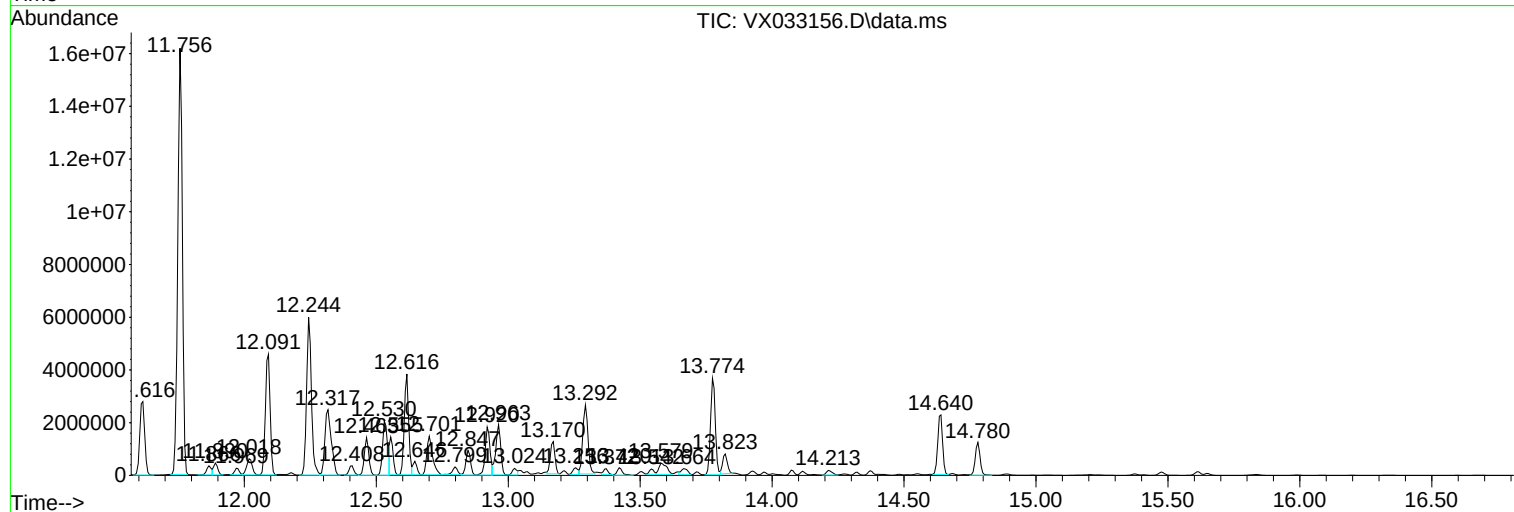
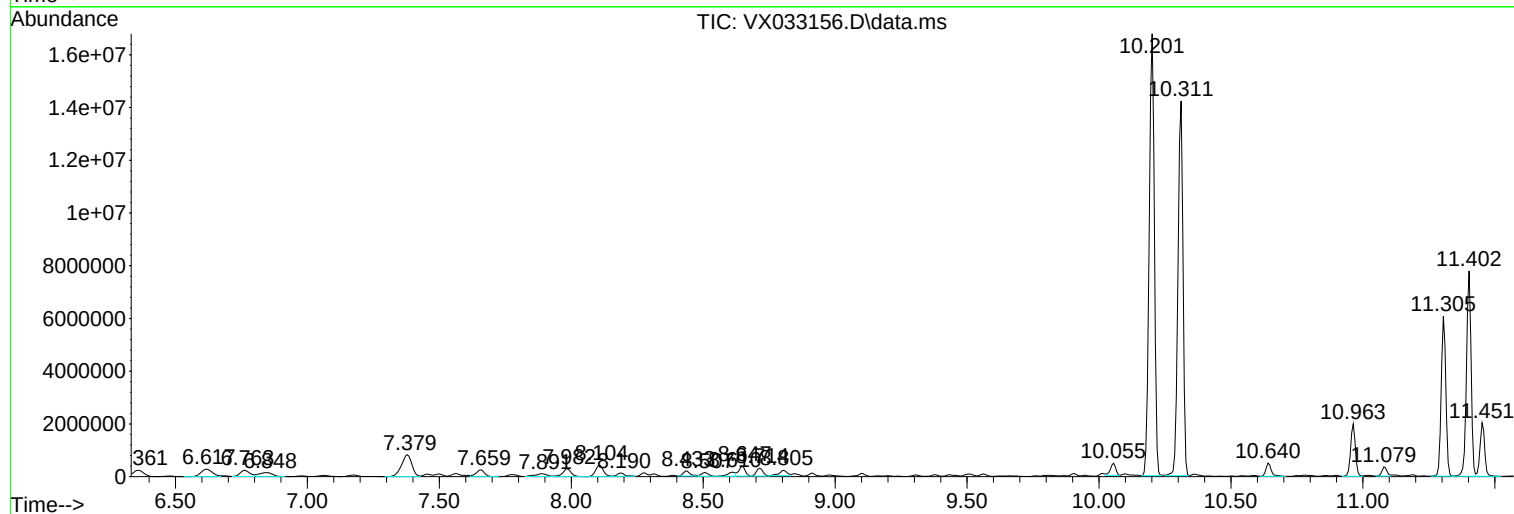
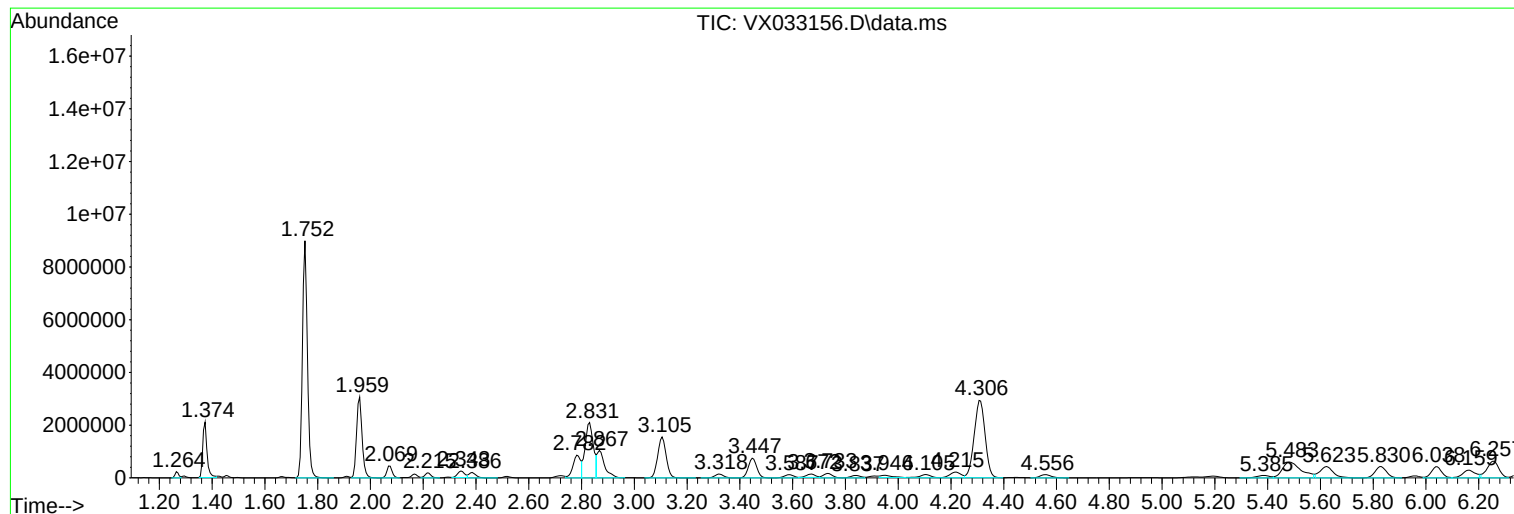
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

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



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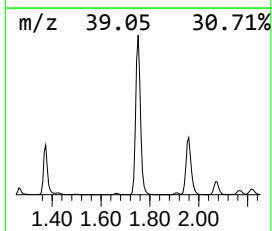
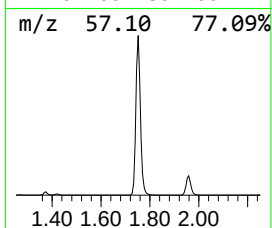
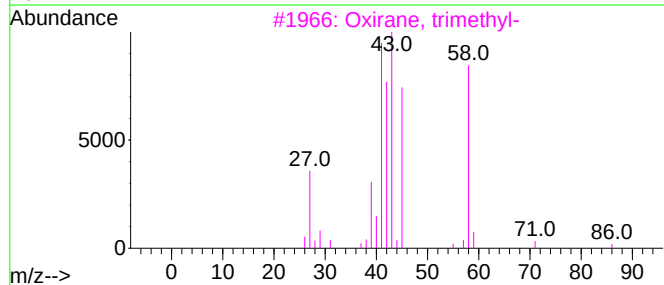
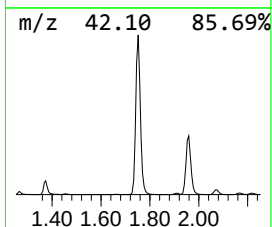
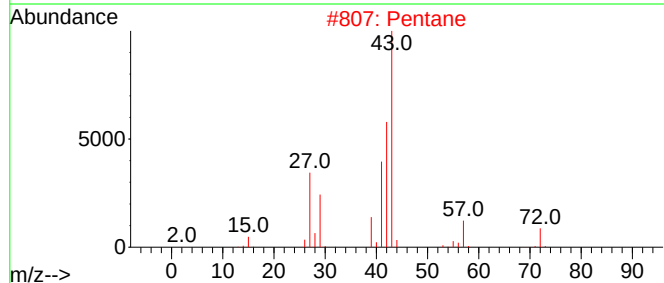
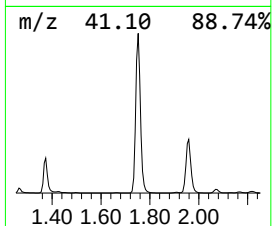
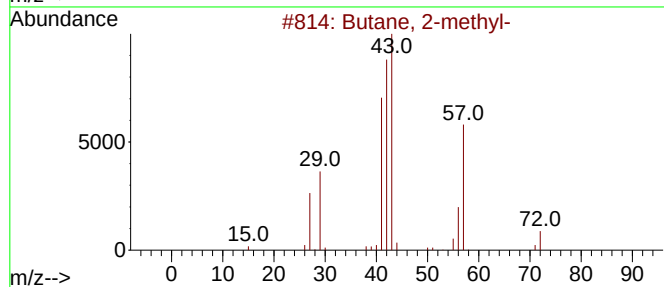
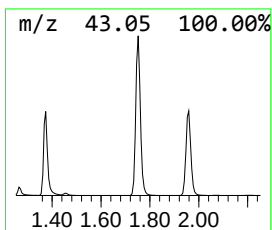
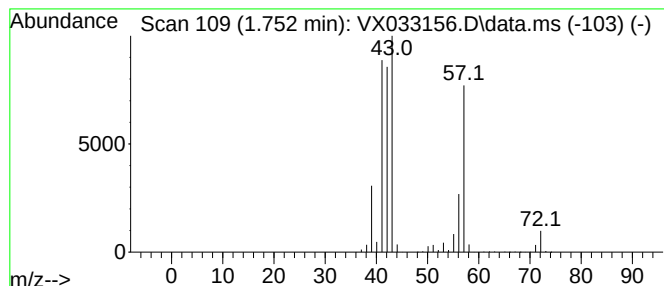
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.752	401.70 ug/l	11828900	Pentafluorobenzene	5.556

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			Oxirane, trimethyl-	86	C5H10O	005076-19-7	33
4			3-Buten-1-ol	72	C4H8O	000627-27-0	25
5			Propane, 2-methyl-2-nitro-	103	C4H9NO2	000594-70-7	10



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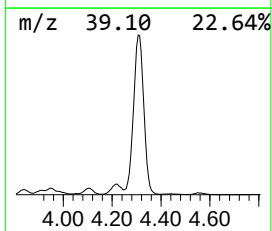
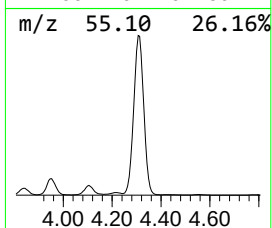
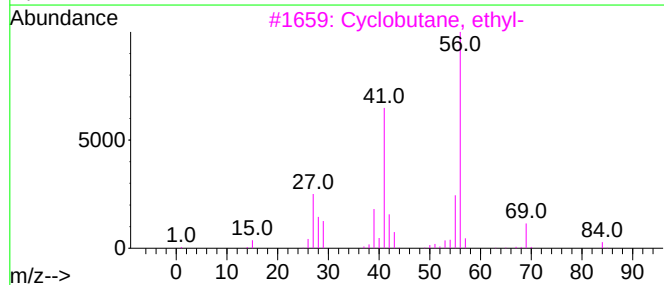
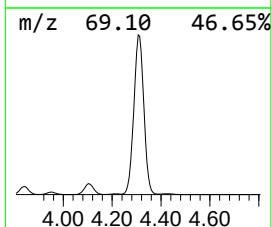
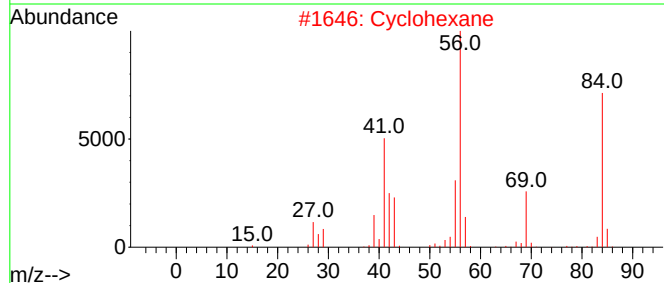
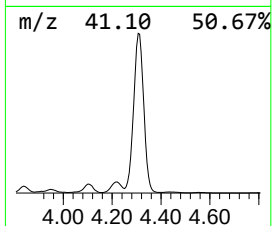
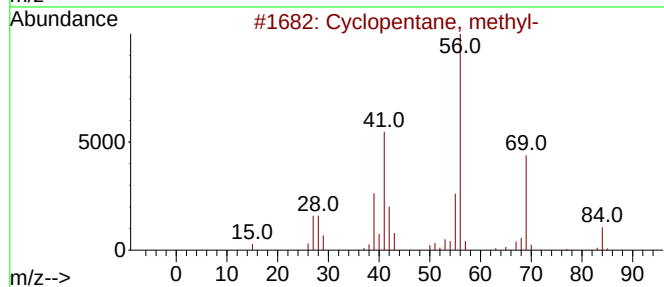
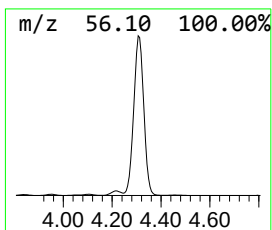
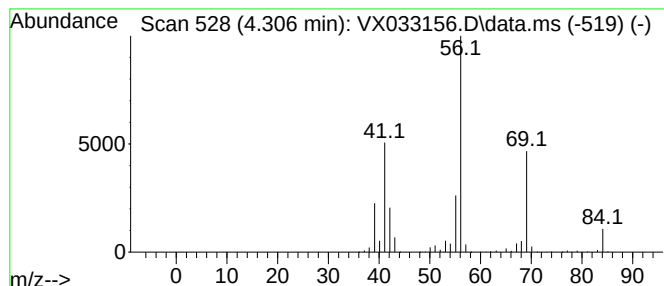
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

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

Peak Number 2 Cyclopentane, methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.306	284.31 ug/l	8372200	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2			Cyclohexane	84	C6H12	000110-82-7	78
3			Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
4			1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	56
5			Cyclobutane	56	C4H8	000287-23-0	53



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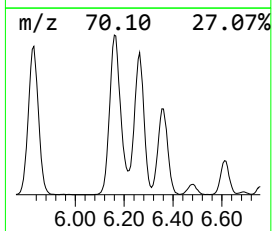
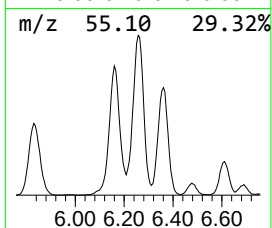
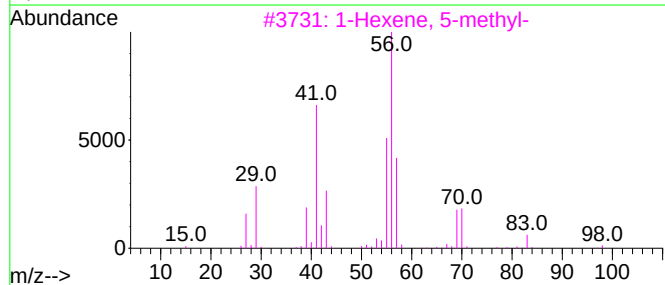
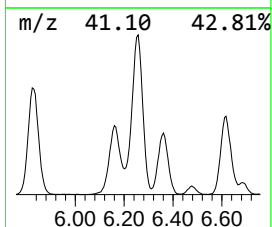
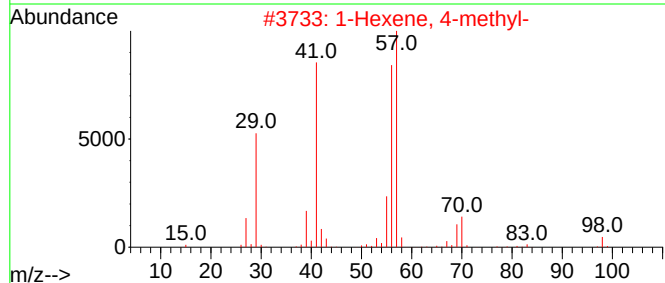
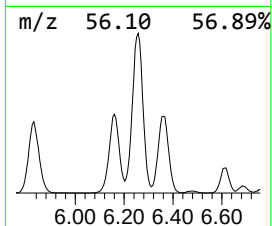
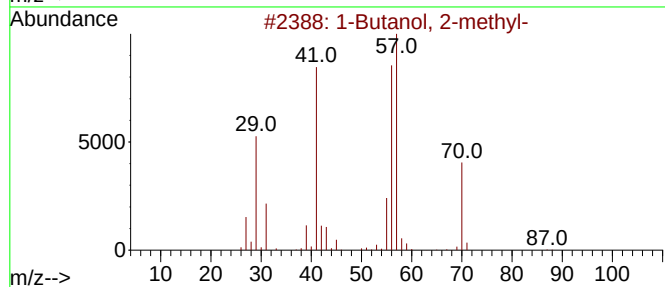
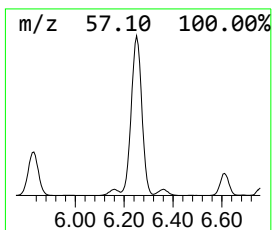
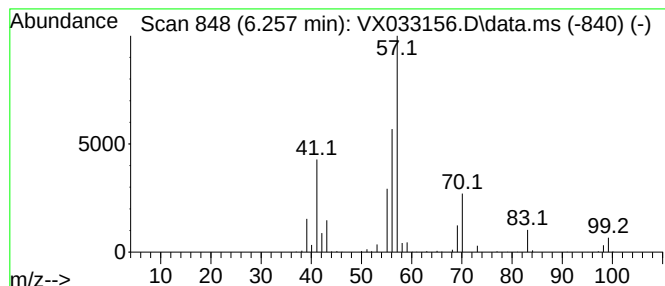
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260

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

Peak Number 3 1-Butanol, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.257	167.28 ug/l	1970160	1,4-Difluorobenzene	6.763

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Butanol, 2-methyl-			88	C5H12O	000137-32-6	59
2	1-Hexene, 4-methyl-			98	C7H14	003769-23-1	43
3	1-Hexene, 5-methyl-			98	C7H14	003524-73-0	40
4	Hexane, 2,2-dimethyl-			114	C8H18	000590-73-8	38
5	Cyclopentane, 1,3-dimethyl-, trans-			98	C7H14	001759-58-6	37



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ClientSampleId :
MW-15

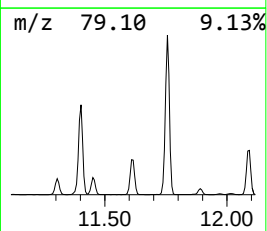
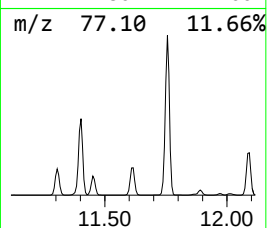
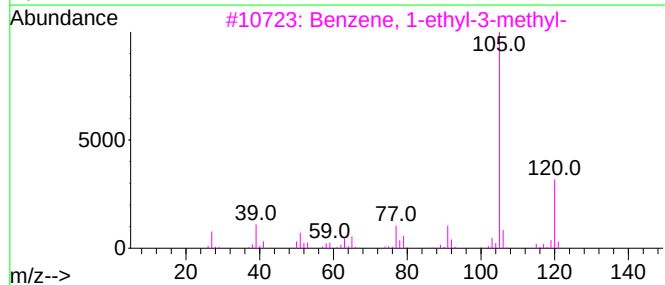
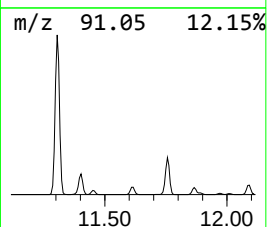
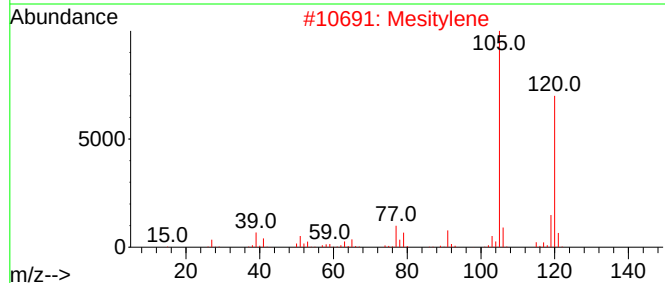
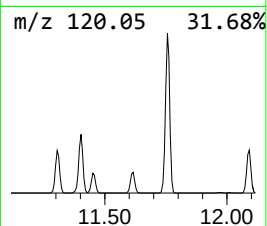
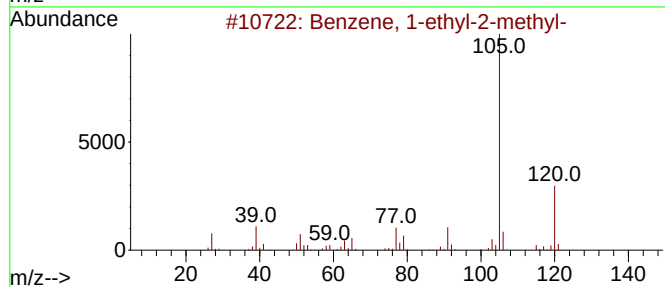
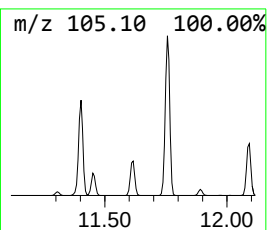
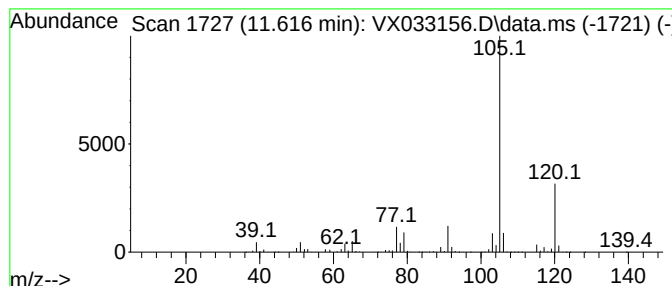
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.616	198.03 ug/l	3579860	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			Mesitylene	120	C9H12	000108-67-8	91
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120122\
Data File : VX033156.D
Acq On : 01 Dec 2022 12:28
Operator : JC/MD
Sample : N5815-25
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-15

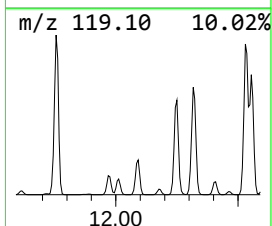
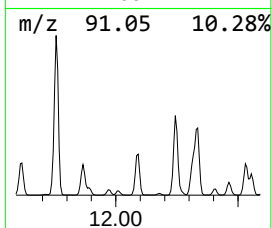
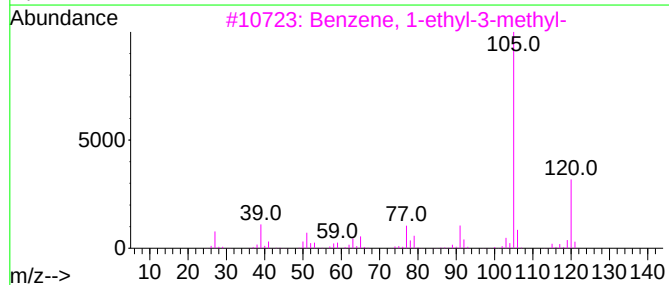
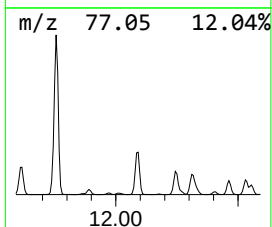
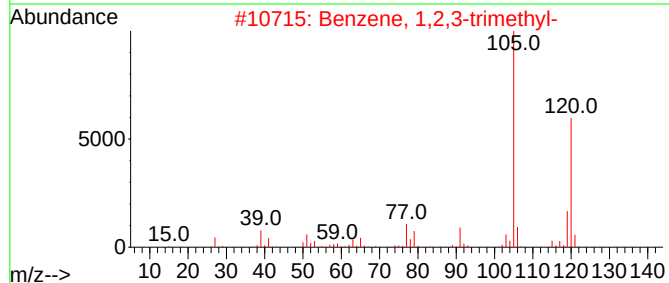
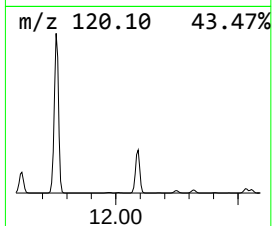
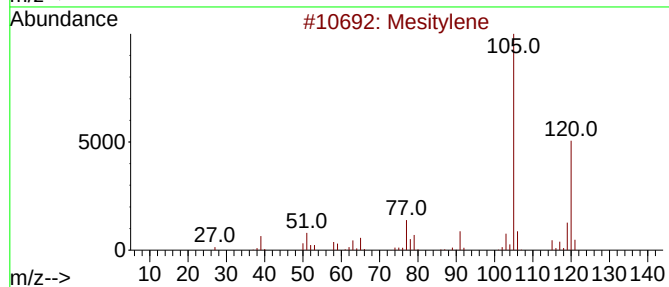
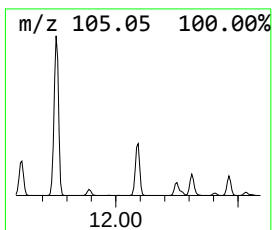
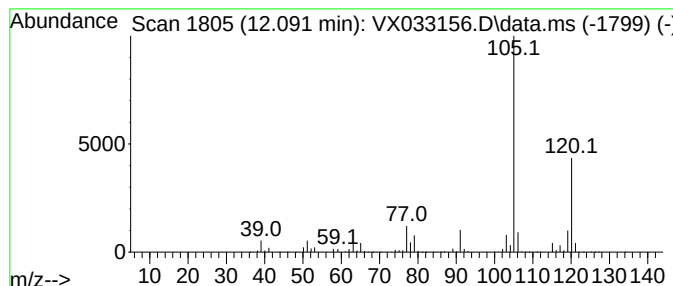
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1,2,3-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.091	316.68 ug/l	5724630	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 : VX033156.D
Acq On : 01 Dec 2022 12:28
Operator : JC/MD
Sample : N5815-25
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-15

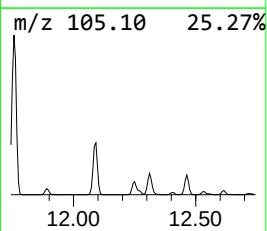
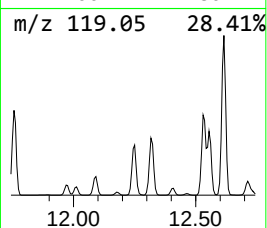
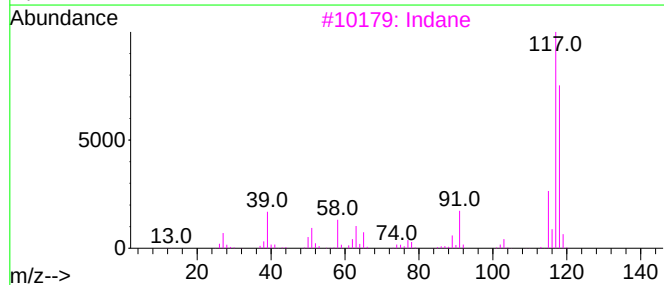
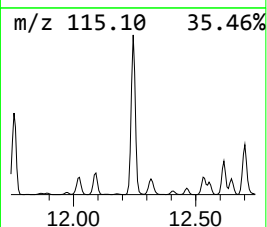
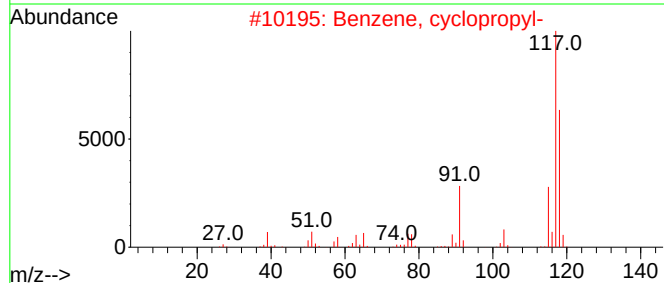
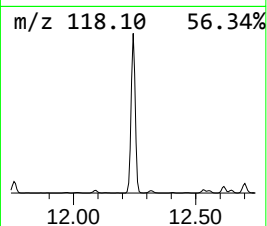
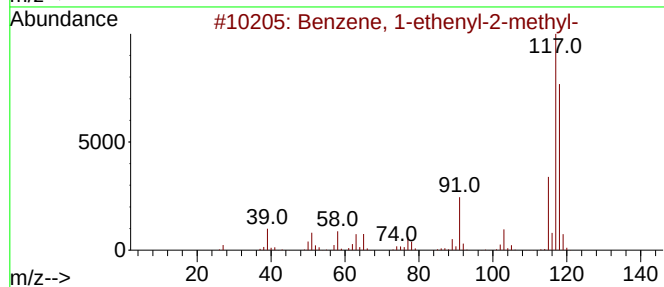
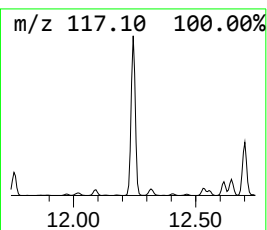
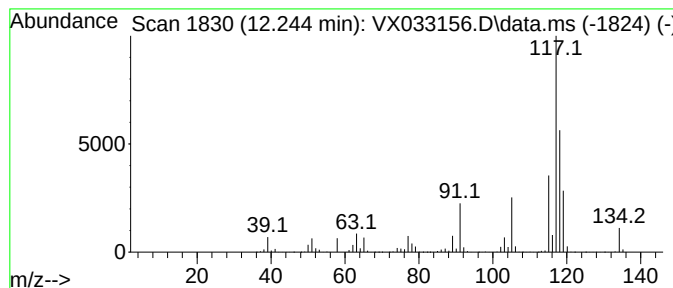
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethenyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	421.78 ug/l	7624610	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	91
2			Benzene, cyclopropyl-	118	C9H10	000873-49-4	91
3			Indane	118	C9H10	000496-11-7	64
4			Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
5			(Z)-1-Phenylpropene	118	C9H10	000766-90-5	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120122\
Data File : VX033156.D
Acq On : 01 Dec 2022 12:28
Operator : JC/MD
Sample : N5815-25
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-15

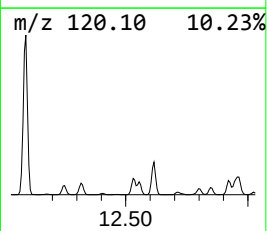
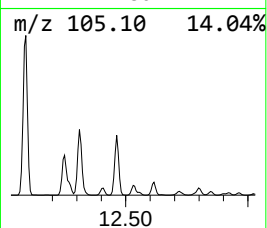
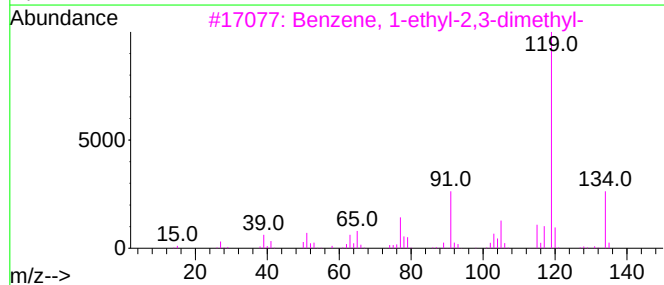
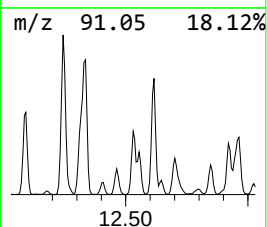
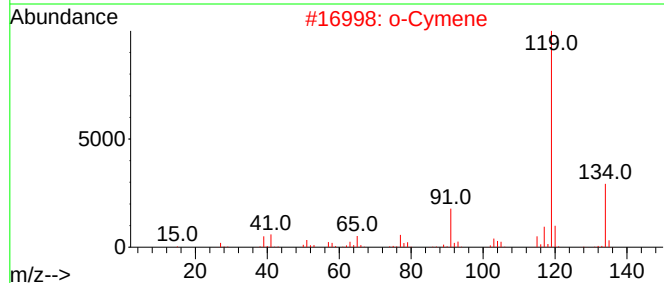
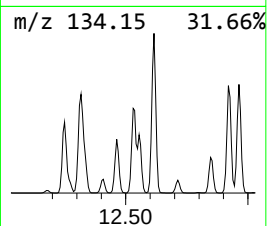
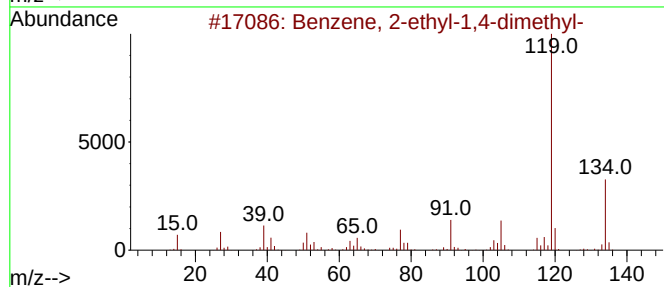
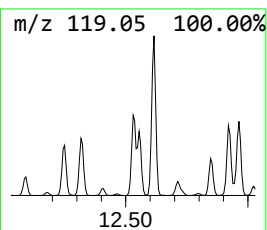
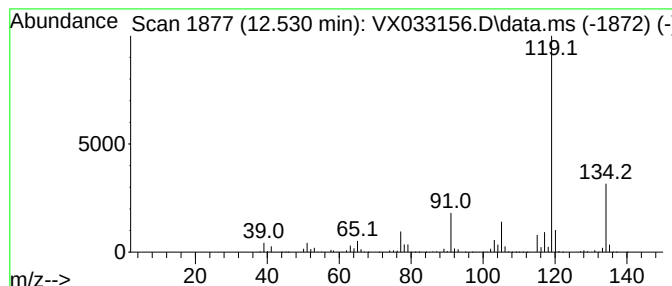
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.530	158.81 ug/l	2870880	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2			o-Cymene	134	C10H14	000527-84-4	95
3			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
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	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120122\
Data File : VX033156.D
Acq On : 01 Dec 2022 12:28
Operator : JC/MD
Sample : N5815-25
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-15

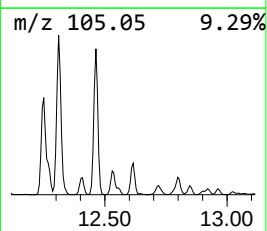
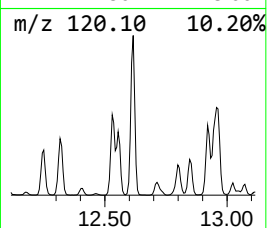
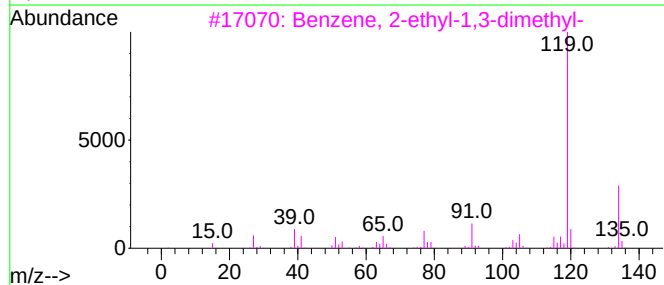
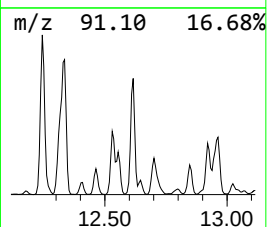
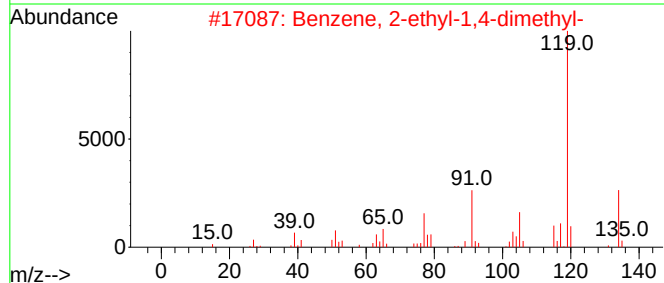
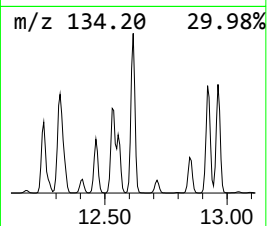
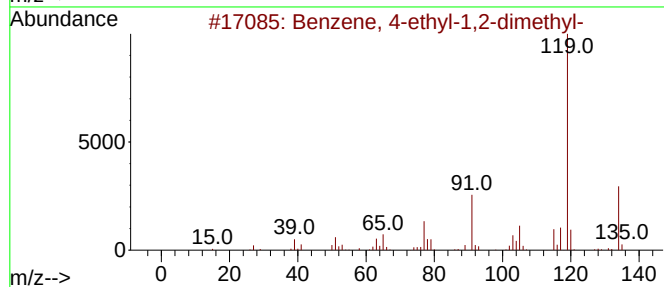
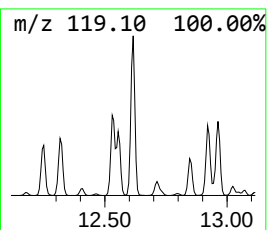
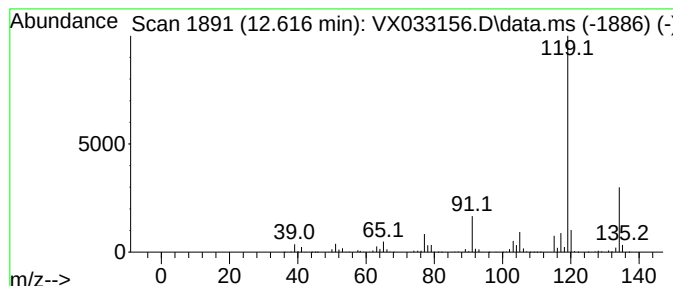
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.616	245.72 ug/l	4441840	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	96
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5			o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120122\
Data File : VX033156.D
Acq On : 01 Dec 2022 12:28
Operator : JC/MD
Sample : N5815-25
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-15

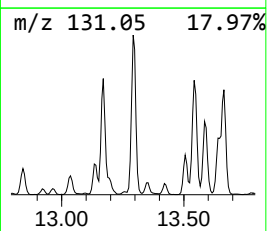
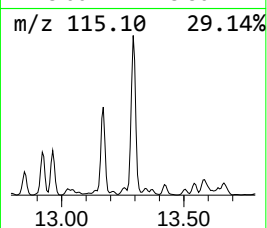
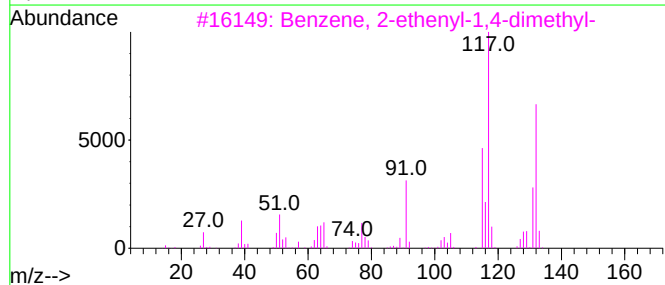
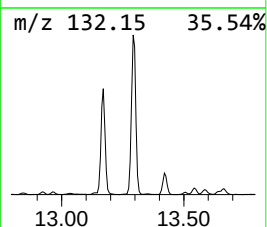
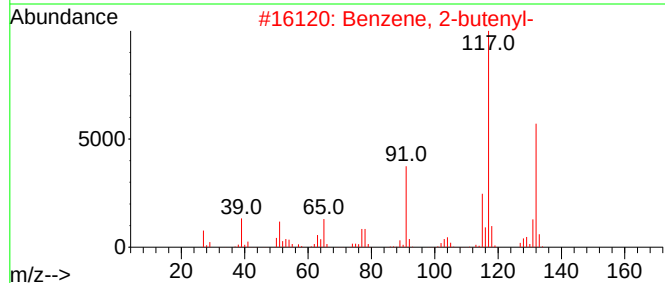
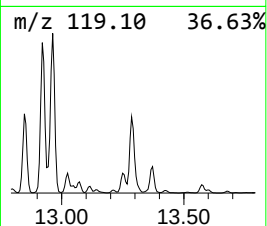
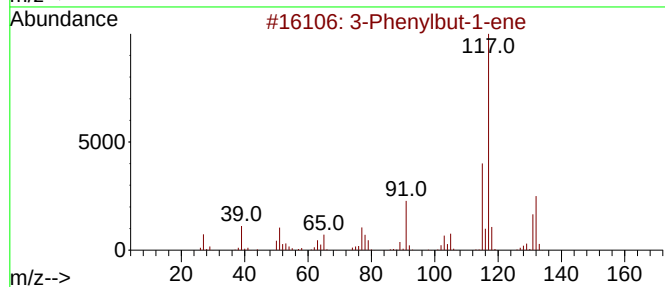
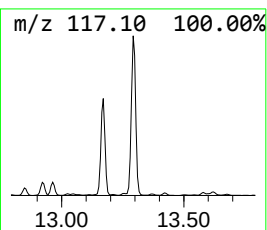
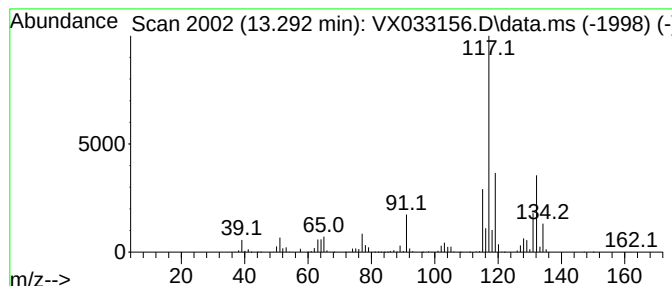
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Quant Title : SW846 8260

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

Peak Number 9 3-Phenylbut-1-ene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	196.25 ug/l	3547620	1,4-Dichlorobenzene-d4	12.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
2		Benzene, 2-butenyl-	132	C10H12	001560-06-1	81
3		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	78
4		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	76
5		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120122\
Data File : VX033156.D
Acq On : 01 Dec 2022 12:28
Operator : JC/MD
Sample : N5815-25
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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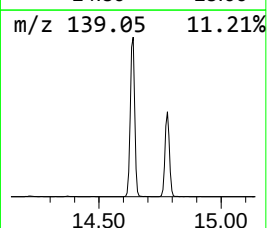
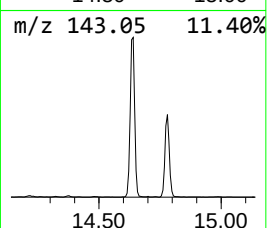
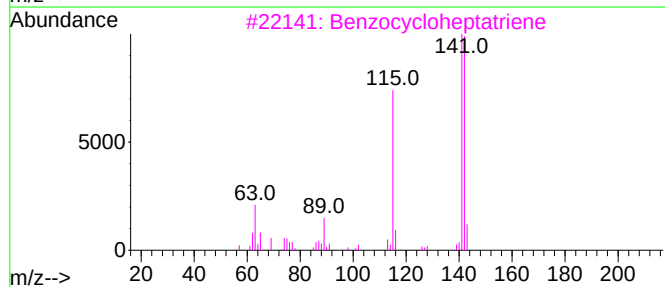
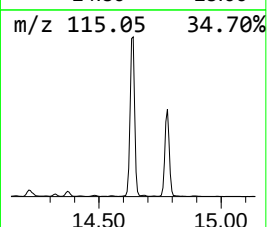
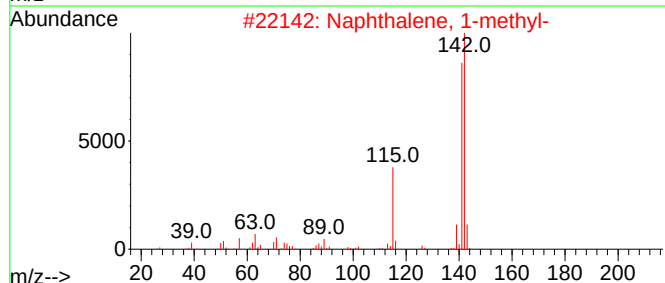
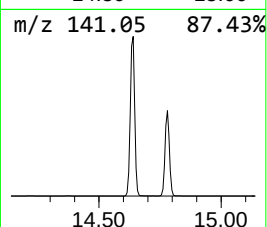
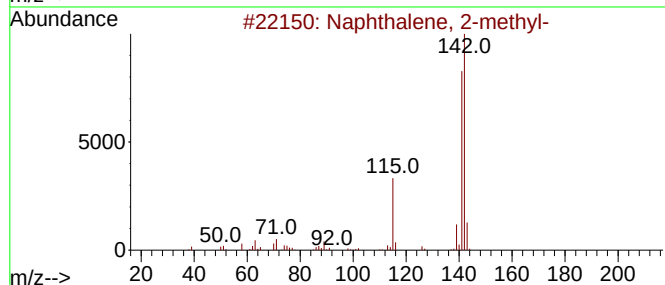
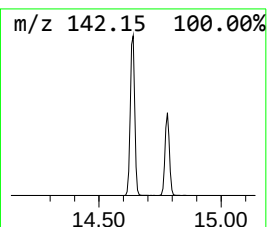
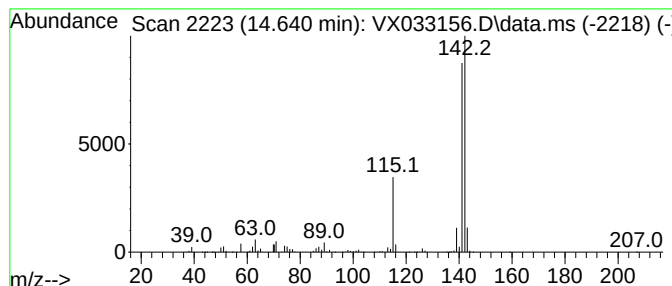
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.640	157.94 ug/l	2855020	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3			Benzocycloheptatriene	142	C11H10	000264-09-5	91
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120122\
Data File : VX033156.D
Acq On : 01 Dec 2022 12:28
Operator : JC/MD
Sample : N5815-25
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-15

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-	1.752	401.7	ug/l	11828900	1	5.556	1472360	50.0
Cyclopentane, m...	4.306	284.3	ug/l	8372200	1	5.556	1472360	50.0
1-Butanol, 2-me...	6.257	167.3	ug/l	1970160	2	6.763	588889	50.0
Benzene, 1-ethy...	11.616	198.0	ug/l	3579860	4	12.024	903858	50.0
Benzene, 1,2,3-...	12.091	316.7	ug/l	5724630	4	12.024	903858	50.0
Benzene, 1-ethe...	12.244	421.8	ug/l	7624610	4	12.024	903858	50.0
Benzene, 2-ethy...	12.530	158.8	ug/l	2870880	4	12.024	903858	50.0
Benzene, 4-ethy...	12.616	245.7	ug/l	4441840	4	12.024	903858	50.0
3-Phenylbut-1-ene	13.292	196.3	ug/l	3547620	4	12.024	903858	50.0
Naphthalene, 2-...	14.640	157.9	ug/l	2855020	4	12.024	903858	50.0