

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061024\
 Data File : VX041811.D
 Acq On : 10 Jun 2024 15:31
 Operator : JC/MD
 Sample : P2767-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-3

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\82X052924W.M
 Title : SW846 8260

Signal : TIC: VX041811.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	53	rBV	73937	76495	1.29%	0.154%
2	1.721	99	104	114	rVB	562980	731651	12.29%	1.478%
3	1.935	135	139	154	rVB	332478	472357	7.94%	0.954%
4	2.057	154	159	165	rBV	52829	73880	1.24%	0.149%
5	2.203	178	183	195	rVB	74710	115109	1.93%	0.232%
6	2.325	195	203	210	rBV	42151	83547	1.40%	0.169%
7	2.709	257	266	270	rBV4	25274	67187	1.13%	0.136%
8	2.770	270	276	279	rVV	121321	254685	4.28%	0.514%
9	2.813	279	283	288	rVV	283969	610664	10.26%	1.233%
10	2.855	288	290	314	rVB2	126298	306101	5.14%	0.618%
11	3.093	319	329	347	rBV	276105	621440	10.44%	1.255%
12	3.319	356	366	377	rBV2	40600	104237	1.75%	0.211%
13	3.434	377	385	400	rVB	109041	244096	4.10%	0.493%
14	3.660	414	422	428	rBV	43812	101961	1.71%	0.206%
15	3.831	441	450	456	rBV2	48814	126333	2.12%	0.255%
16	4.093	486	493	504	rVB3	32203	82382	1.38%	0.166%
17	4.294	516	526	539	rVV	533900	1520964	25.55%	3.072%
18	4.422	540	547	561	rVB5	19297	62317	1.05%	0.126%
19	5.190	664	673	689	rVB4	35603	127231	2.14%	0.257%
20	5.385	689	705	710	rBV2	84293	261215	4.39%	0.528%
21	5.465	710	718	727	rVV	318341	940978	15.81%	1.900%
22	5.550	727	732	739	rVB	79806	192826	3.24%	0.389%
23	5.830	765	778	790	rBV4	78618	273211	4.59%	0.552%
24	5.958	790	799	813	rVB	81321	215315	3.62%	0.435%
25	6.153	813	831	841	rBV3	96600	382638	6.43%	0.773%
26	6.257	841	848	855	rVV3	91322	261021	4.39%	0.527%
27	6.355	855	864	875	rVB	115631	329466	5.54%	0.665%
28	6.757	923	930	938	rBV	198096	516129	8.67%	1.042%
29	6.836	940	943	953	rVB5	57331	150018	2.52%	0.303%
30	7.172	989	998	1018	rVB4	46251	118686	1.99%	0.240%
31	7.379	1018	1032	1046	rBV	600402	1485639	24.96%	3.000%
32	7.653	1070	1077	1089	rVB	66826	134334	2.26%	0.271%
33	7.879	1103	1114	1121	rBV2	50406	166646	2.80%	0.337%
34	8.141	1152	1157	1163	rVB	53550	88240	1.48%	0.178%
35	8.501	1210	1216	1225	rVB2	111053	205083	3.45%	0.414%
36	8.598	1225	1232	1235	rBV3	53666	104672	1.76%	0.211%

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Integrator: RTE

Smoothing : ON

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Sampling : 1

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Stop Thrs : 0

Peak Location: TOP

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37	8.647	1235	1240	1247	rVB	439309	688048	11.56%	1.390%
38	8.720	1247	1252	1257	rBV	229143	350127	5.88%	0.707%
39	8.970	1289	1293	1300	rVB	48261	78746	1.32%	0.159%
40	9.098	1306	1314	1319	rVV3	36530	69080	1.16%	0.140%
41	9.159	1319	1324	1335	rVB	72810	124944	2.10%	0.252%
42	10.055	1461	1471	1476	rBV	460666	653478	10.98%	1.320%
43	10.122	1476	1482	1485	rVV	33152	60269	1.01%	0.122%
44	10.195	1489	1494	1499	rVB	775033	1033651	17.37%	2.087%
45	10.299	1505	1511	1527	rVB	4316215	5952284	100.00%	12.021%
46	10.640	1562	1567	1578	rVB	2340311	2944907	49.48%	5.947%
47	10.964	1614	1620	1628	rBV	303904	403097	6.77%	0.814%
48	11.079	1635	1639	1644	rBV	416469	516895	8.68%	1.044%
49	11.305	1667	1676	1683	rBV	526964	651916	10.95%	1.317%
50	11.378	1683	1688	1696	rVV	1850663	3162334	53.13%	6.386%
51	11.451	1696	1700	1707	rVB	784043	949754	15.96%	1.918%
52	11.610	1721	1726	1737	rBV	1356549	1680512	28.23%	3.394%
53	11.750	1744	1749	1756	rBV	3870543	4571523	76.80%	9.232%
54	11.866	1762	1768	1770	rBV	48639	70460	1.18%	0.142%
55	11.890	1770	1772	1780	rVB	71271	84476	1.42%	0.171%
56	11.969	1780	1785	1788	rBV	159916	189633	3.19%	0.383%
57	12.018	1788	1793	1799	rVB2	547499	744220	12.50%	1.503%
58	12.085	1799	1804	1809	rBV	1060403	1266471	21.28%	2.558%
59	12.244	1824	1830	1832	rBV2	611644	803081	13.49%	1.622%
60	12.268	1832	1834	1838	rVV	433185	466206	7.83%	0.941%
61	12.317	1838	1842	1850	rVB3	428107	677465	11.38%	1.368%
62	12.408	1851	1857	1861	rVB2	72396	101678	1.71%	0.205%
63	12.463	1861	1866	1872	rVB	191511	228351	3.84%	0.461%
64	12.530	1872	1877	1879	rBV	435989	528611	8.88%	1.068%
65	12.555	1879	1881	1885	rVB	284715	294435	4.95%	0.595%
66	12.610	1885	1890	1894	rBV	776495	964271	16.20%	1.947%
67	12.646	1894	1896	1900	rVB	105886	109307	1.84%	0.221%
68	12.695	1900	1904	1912	rBV2	640996	949573	15.95%	1.918%
69	12.847	1924	1929	1934	rVB2	213792	279927	4.70%	0.565%
70	12.920	1934	1941	1945	rBV	483175	587263	9.87%	1.186%
71	12.963	1945	1948	1954	rVB	382639	443098	7.44%	0.895%
72	13.042	1954	1961	1967	rBV4	61580	146510	2.46%	0.296%
73	13.134	1973	1976	1978	rBV	53901	62157	1.04%	0.126%
74	13.170	1978	1982	1986	rVB	351375	438529	7.37%	0.886%
75	13.292	1997	2002	2007	rVV2	827655	1100061	18.48%	2.222%
76	13.420	2019	2023	2031	rVB2	135326	178588	3.00%	0.361%

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Title : SW846 8260

77	13.506	2031	2037	2040	rBV2	76234	108499	1.82%	0.219%
78	13.542	2040	2043	2046	rVV	181891	235808	3.96%	0.476%
79	13.585	2046	2050	2055	rVV3	161235	284928	4.79%	0.575%
80	13.664	2055	2063	2068	rVB2	189271	397755	6.68%	0.803%
81	13.774	2074	2081	2089	rBV	1030914	1273859	21.40%	2.573%
82	13.865	2089	2096	2100	rVV3	46033	79433	1.33%	0.160%
83	13.926	2101	2106	2110	rVV2	50924	69437	1.17%	0.140%
84	13.969	2110	2113	2117	rVB	54800	62868	1.06%	0.127%
85	14.073	2126	2130	2137	rBV	84745	111010	1.86%	0.224%
86	14.213	2149	2153	2159	rBV2	68947	121629	2.04%	0.246%
87	14.371	2175	2179	2184	rVB	51445	61165	1.03%	0.124%
88	14.634	2218	2222	2233	rVB	211066	299978	5.04%	0.606%
89	14.780	2241	2246	2256	rVB	149798	196498	3.30%	0.397%

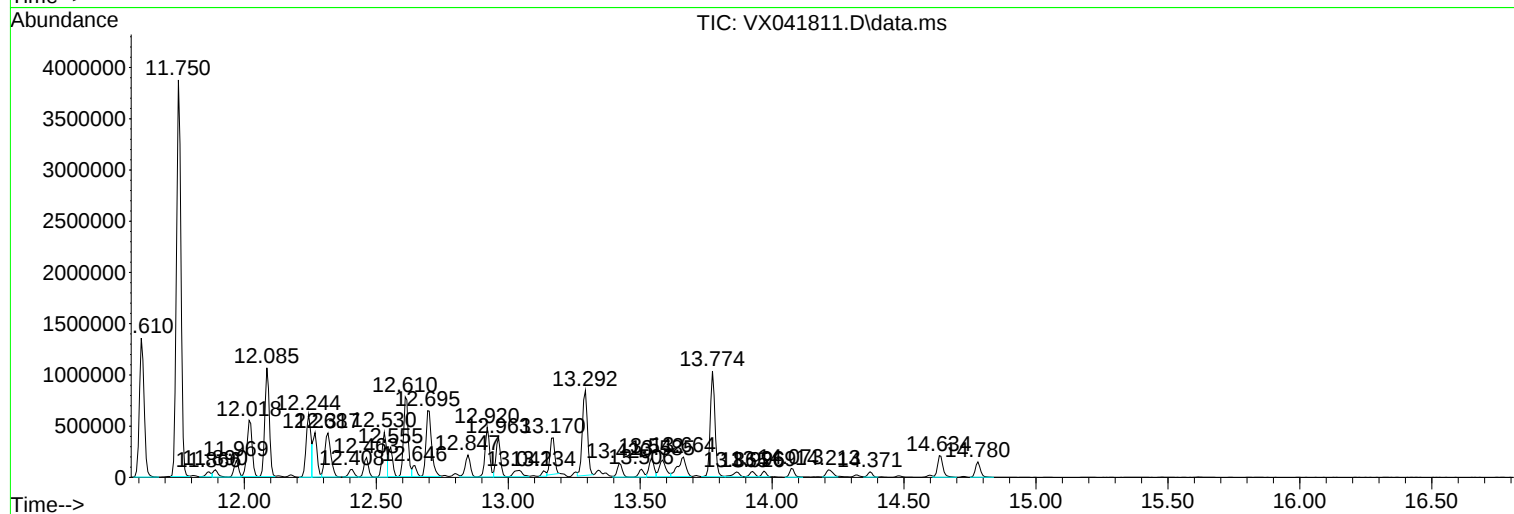
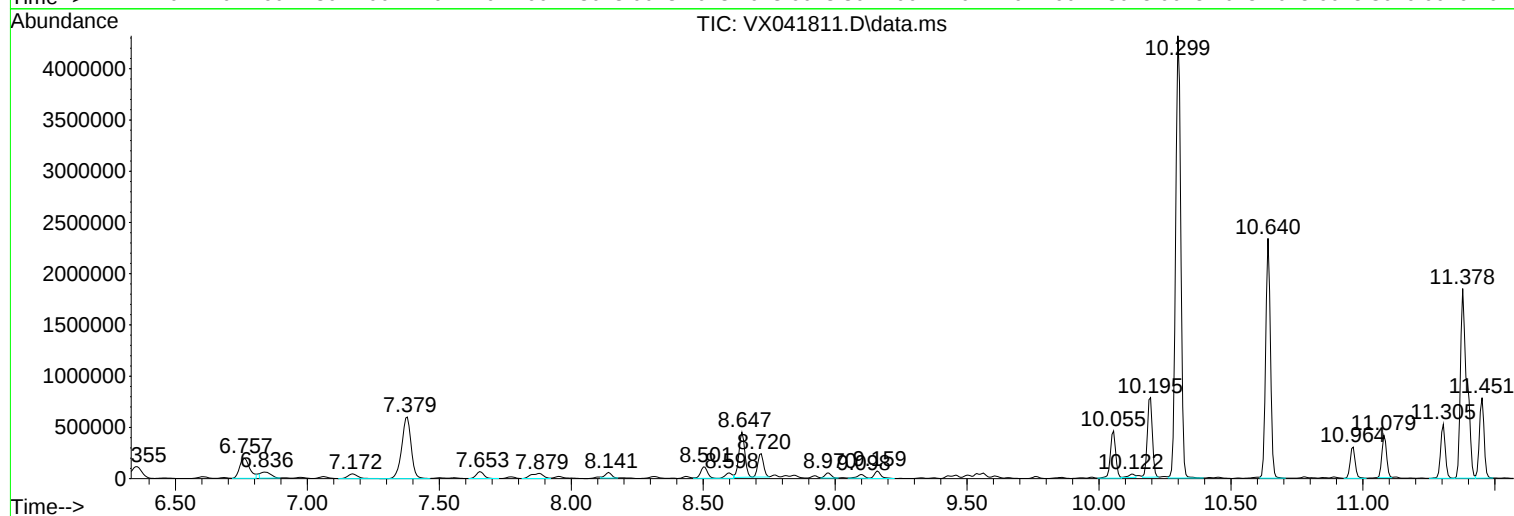
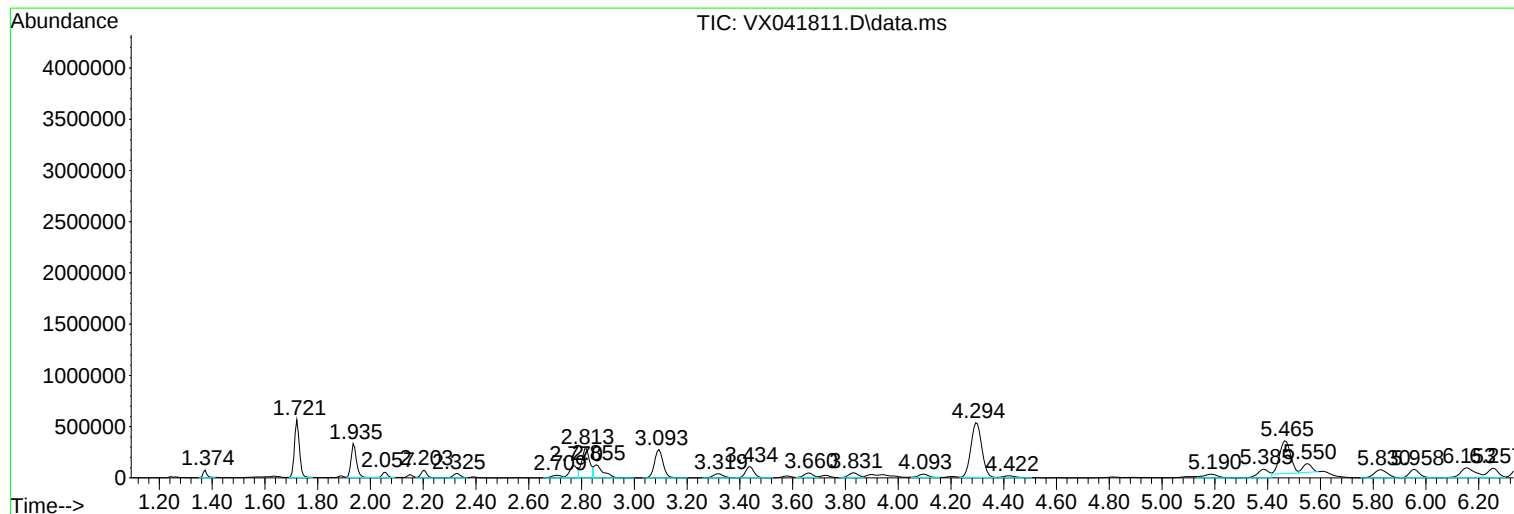
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Instrument :
MSVOA_X
ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X052924W.M
Quant Title : SW846 8260

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



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Instrument :
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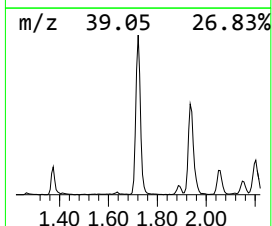
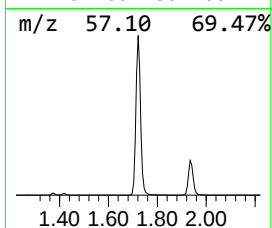
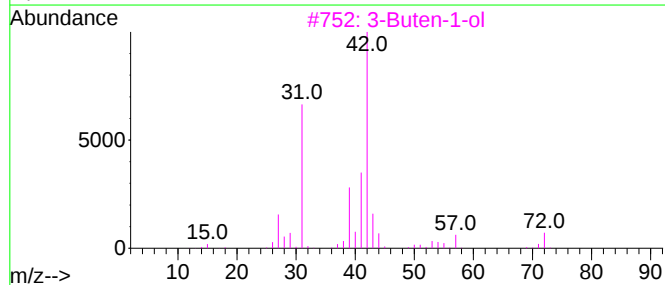
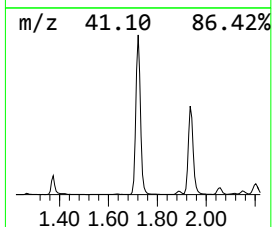
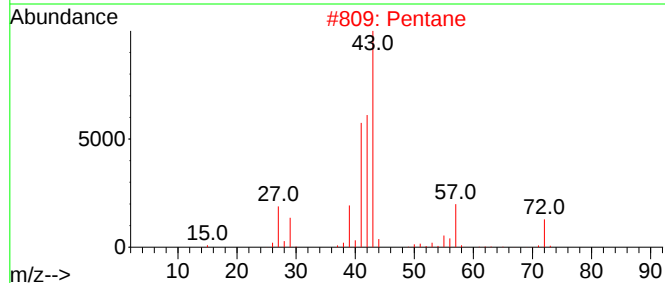
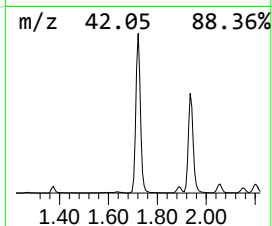
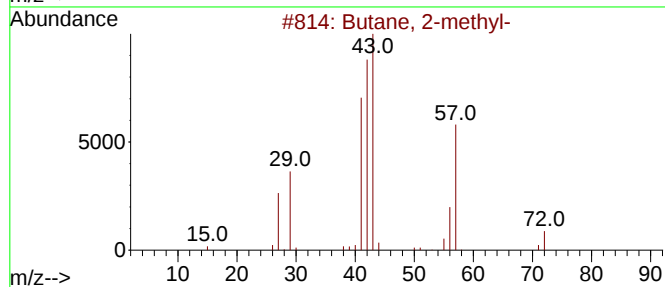
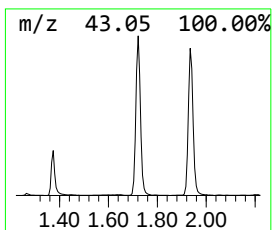
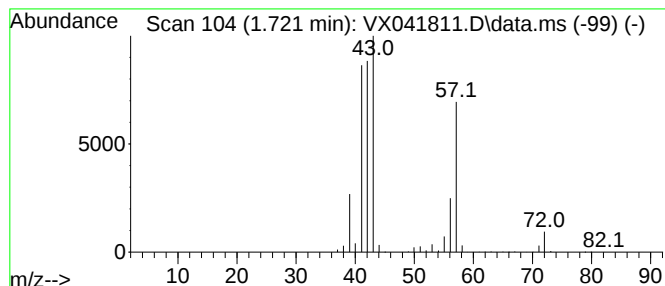
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X052924W.M
Quant Title : SW846 8260

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

Peak Number 1 Butane, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.721	189.72 ug/l	731651	Pentafluorobenzene	5.550

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-			72	C5H12	000078-78-4	91
2	Pentane			72	C5H12	000109-66-0	43
3	3-Buten-1-ol			72	C4H8O	000627-27-0	28
4	1-Butene			56	C4H8	000106-98-9	10
5	Methane, isocyanato-			57	C2H3NO	000624-83-9	9



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ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
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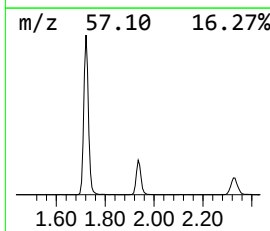
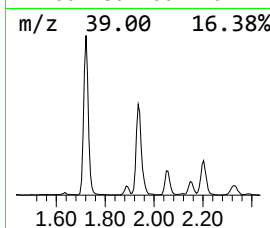
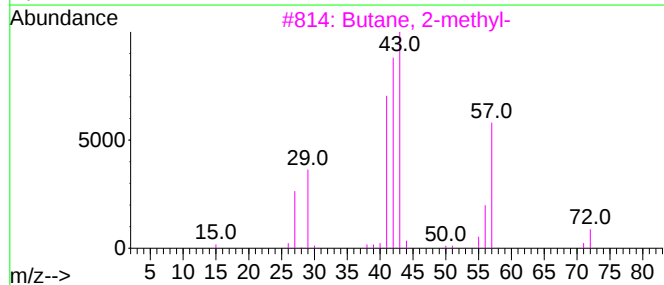
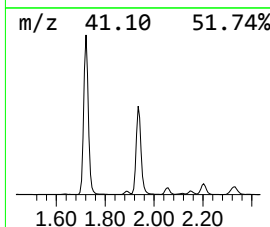
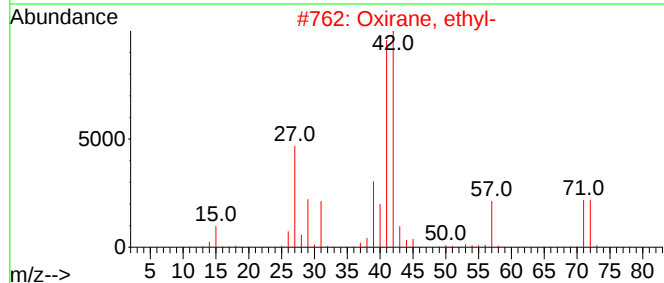
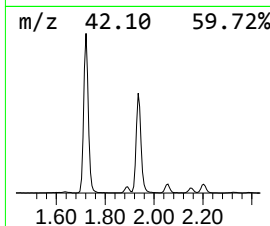
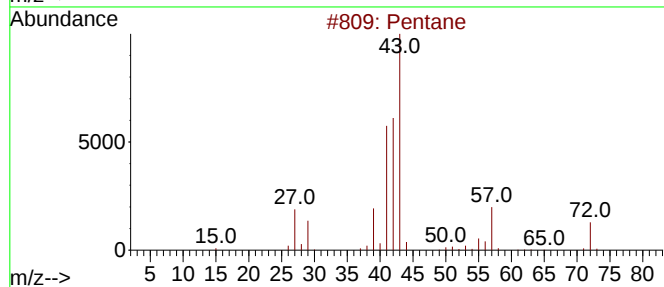
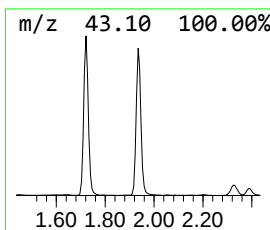
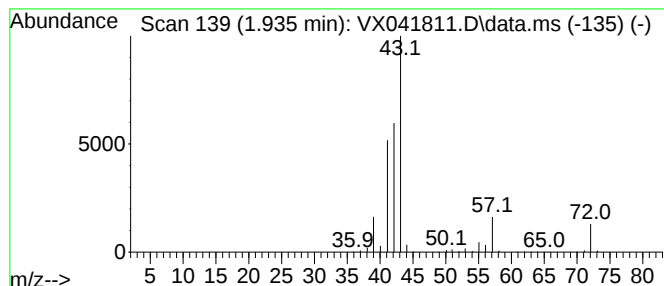
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X052924W.M
Quant Title : SW846 8260

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

Peak Number 2 Pentane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.935	122.48 ug/l	472357	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	39
4			Diaziridine,3,3-dimethyl-	72	C3H8N2	004901-76-2	32
5			2-Propanone, hydrazone	72	C3H8N2	005281-20-9	9



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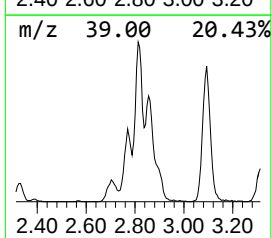
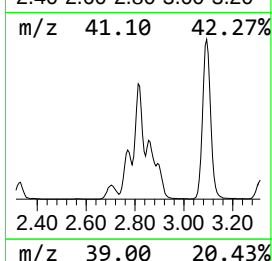
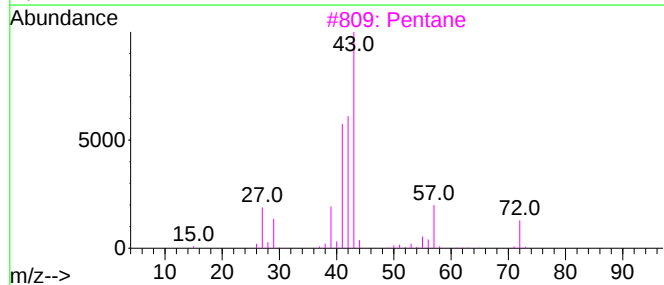
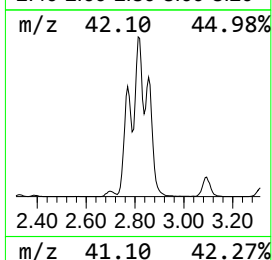
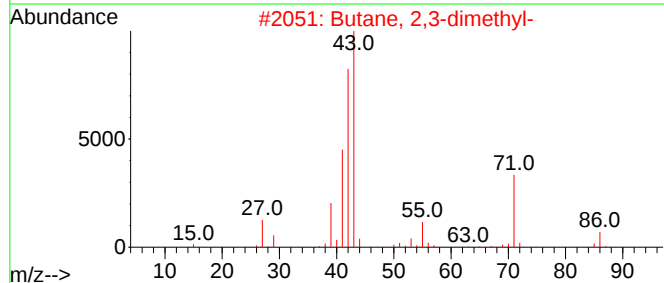
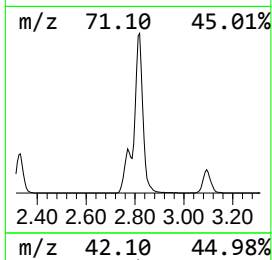
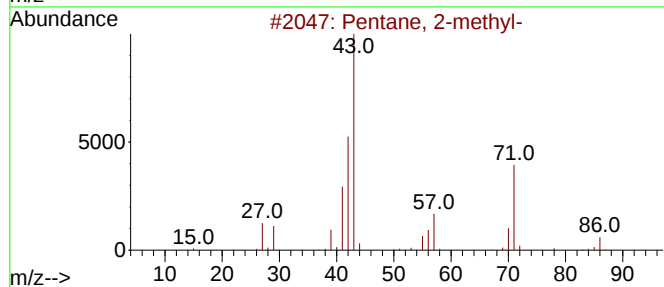
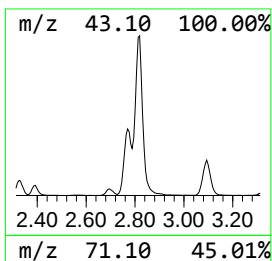
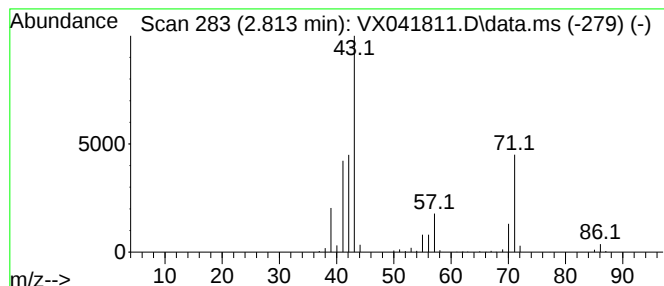
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X052924W.M
Quant Title : SW846 8260

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

Peak Number 3 Pentane, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.813	158.35 ug/l	610664	Pentafluorobenzene	5.550

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2-methyl-	86	C6H14	000107-83-5	91
2			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	47
3			Pentane	72	C5H12	000109-66-0	46
4			1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	40
5			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	38



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ClientSampleId :
MW-3

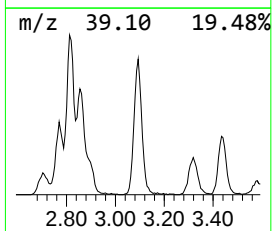
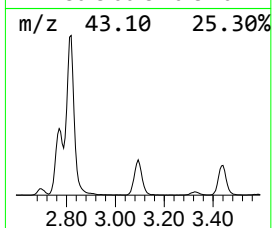
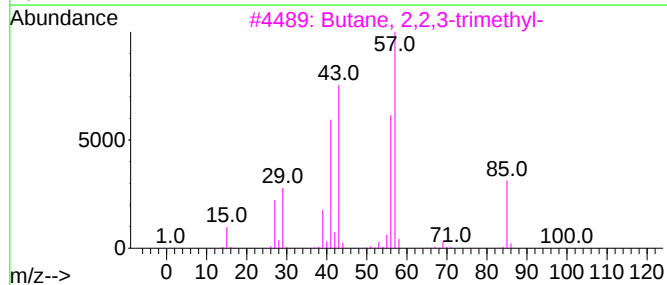
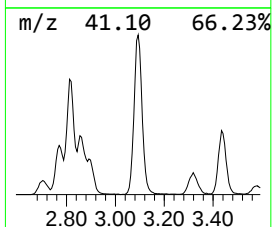
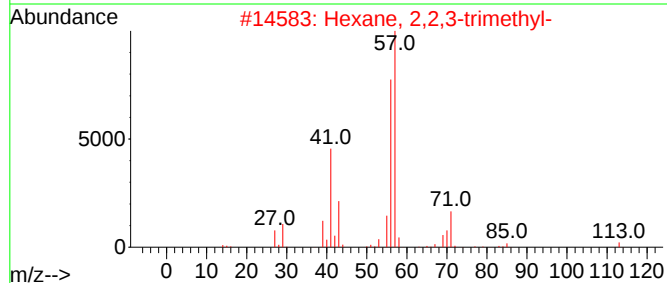
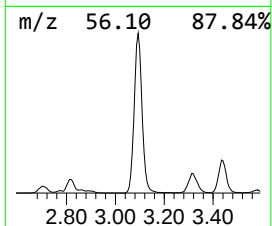
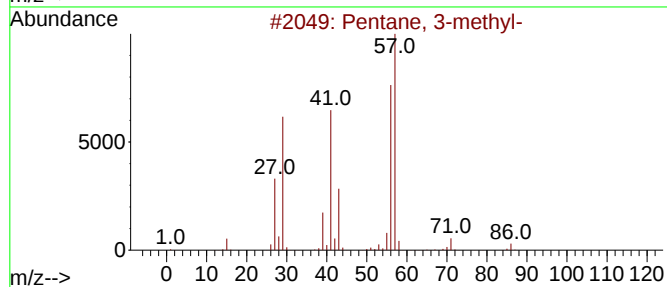
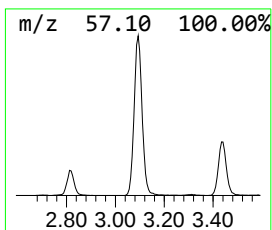
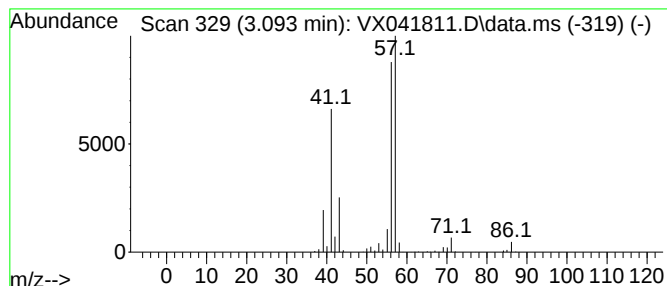
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Quant Title : SW846 8260

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

Peak Number 4 Pentane, 3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.093	161.14 ug/l	621440	Pentafluorobenzene	5.550

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2			Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	78
3			Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	64
4			Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	64
5			3,4-Dimethyldihydrofuran-2,5-dione	128	C6H8O3	007475-92-5	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061024\
Data File : VX041811.D
Acq On : 10 Jun 2024 15:31
Operator : JC/MD
Sample : P2767-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

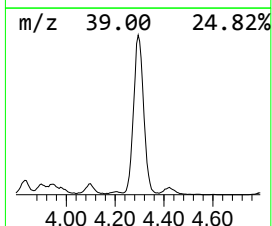
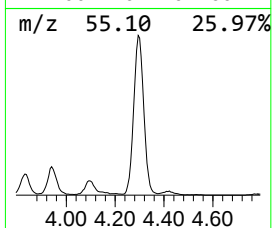
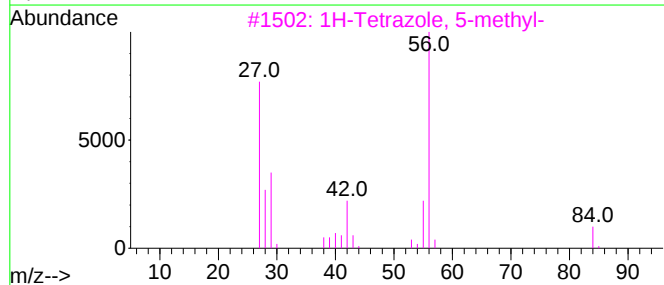
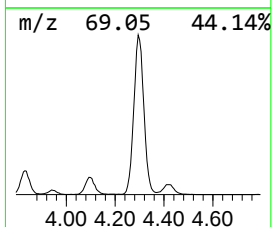
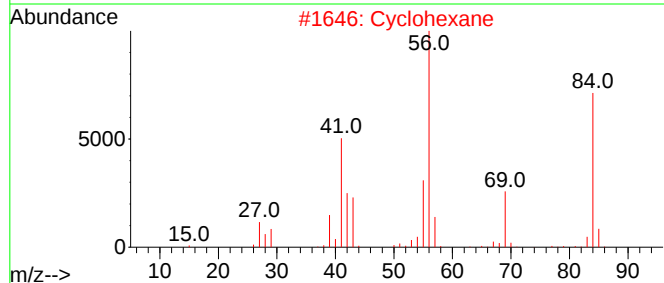
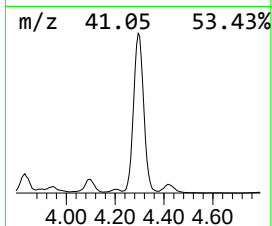
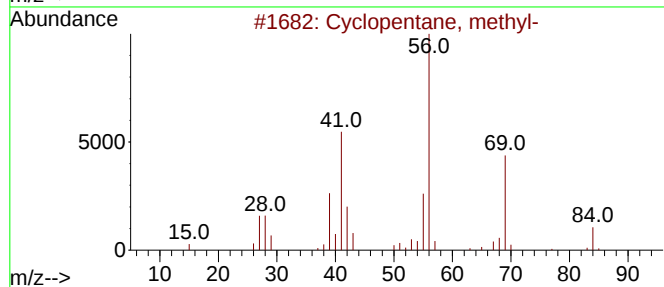
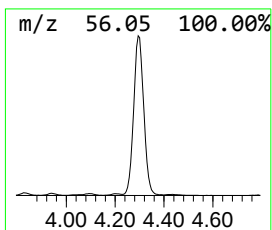
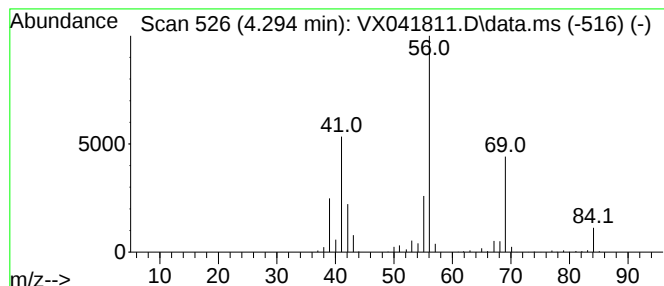
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Quant Title : SW846 8260

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

Peak Number 5 Cyclopentane, methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.294	394.39 ug/l	1520960	Pentafluorobenzene	5.550

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			1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
4			Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
5			Cyclobutane	56	C4H8	000287-23-0	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061024\
Data File : VX041811.D
Acq On : 10 Jun 2024 15:31
Operator : JC/MD
Sample : P2767-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

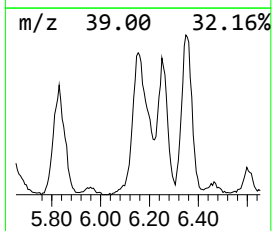
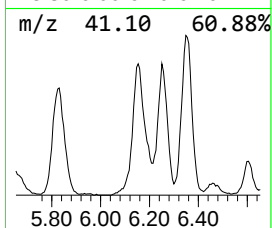
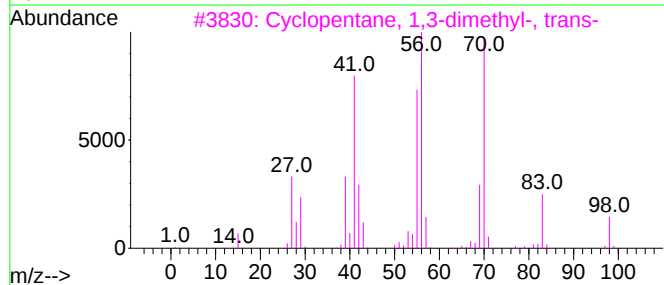
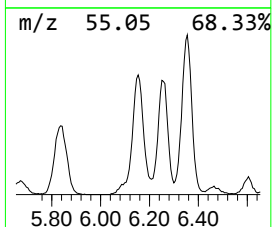
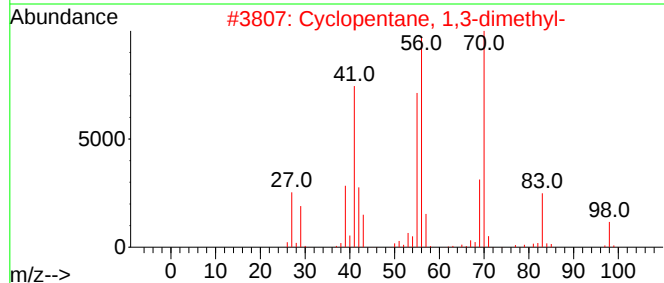
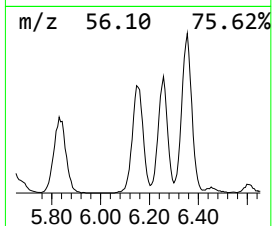
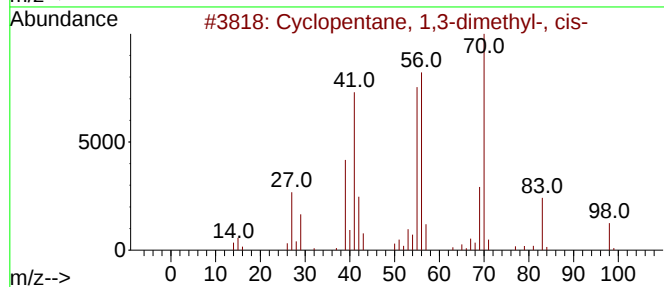
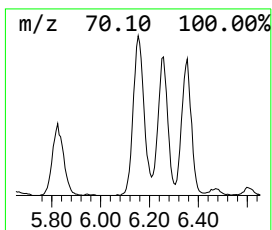
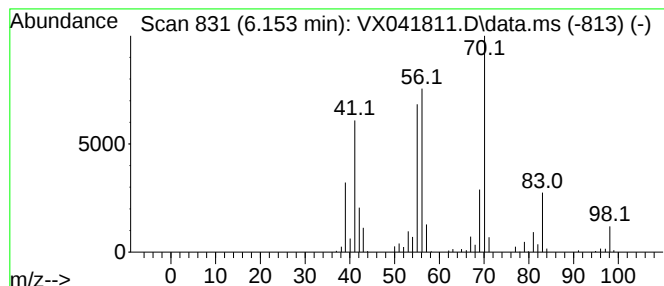
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Quant Title : SW846 8260

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

Peak Number 6 Cyclopentane, 1,3-dimethyl-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.153	99.22 ug/l	382638	Pentafluorobenzene	5.550

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	96
2		Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	95
3		Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	94
4		Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	76
5		Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061024\
Data File : VX041811.D
Acq On : 10 Jun 2024 15:31
Operator : JC/MD
Sample : P2767-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

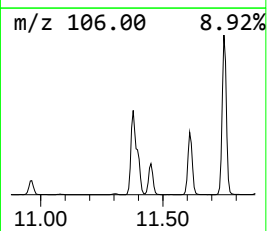
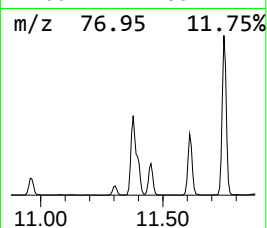
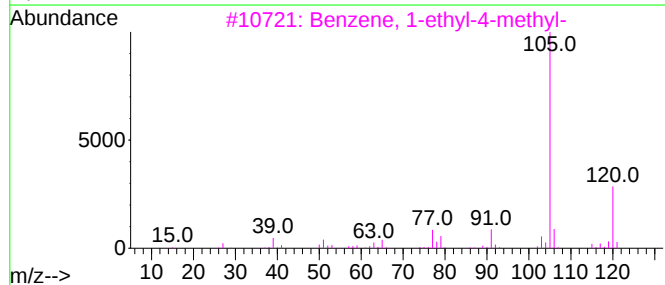
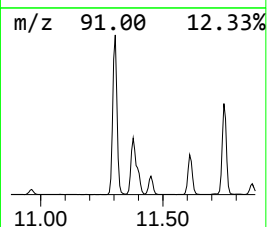
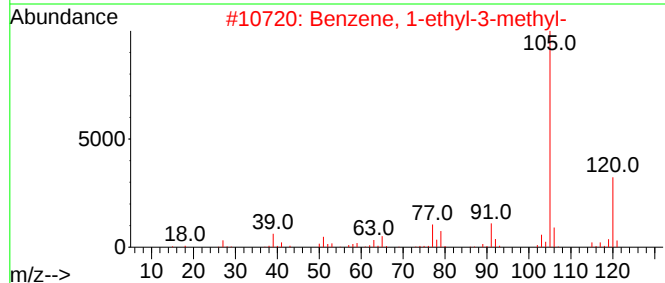
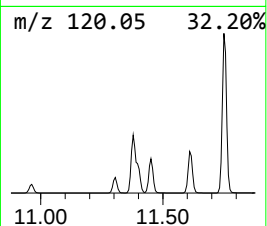
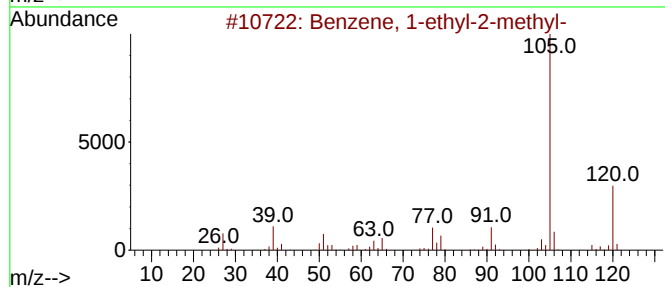
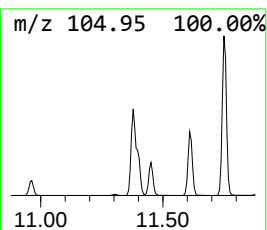
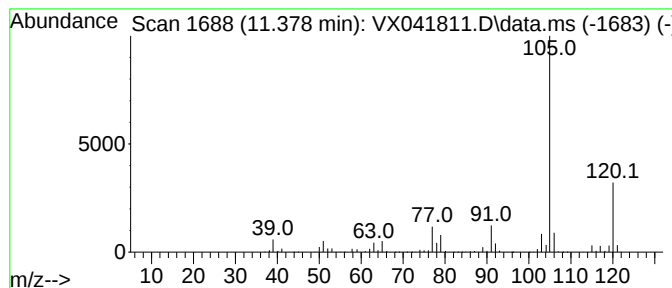
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.378	212.46 ug/l	3162330	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	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\VX061024\
Data File : VX041811.D
Acq On : 10 Jun 2024 15:31
Operator : JC/MD
Sample : P2767-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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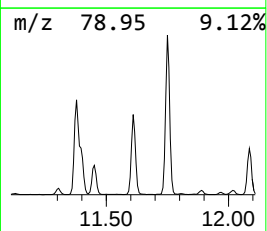
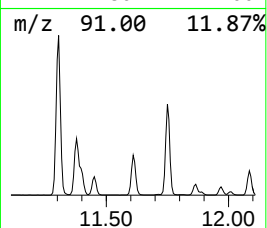
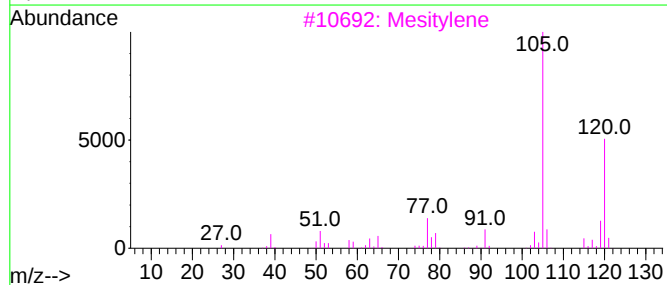
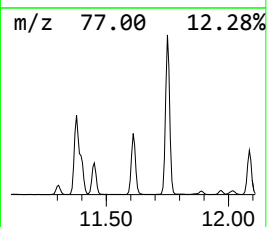
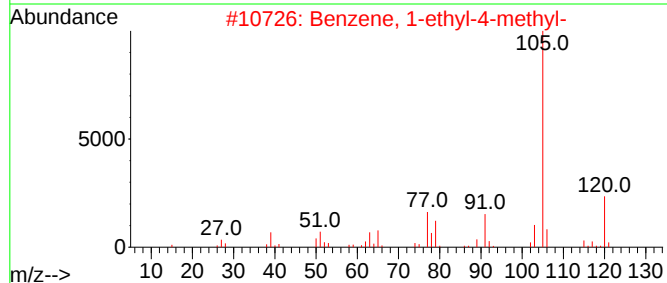
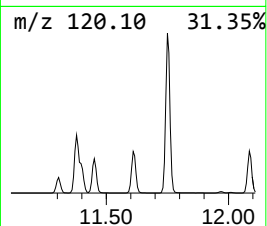
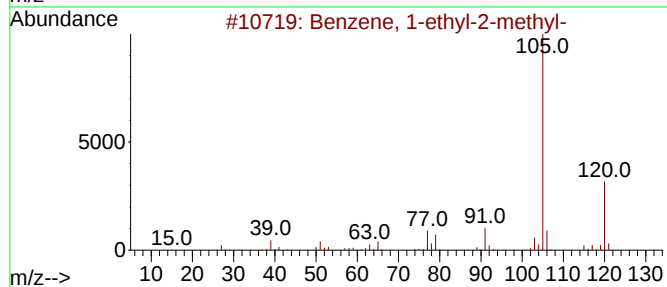
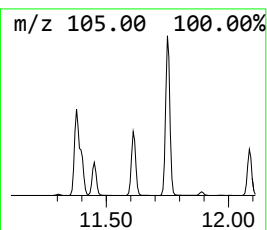
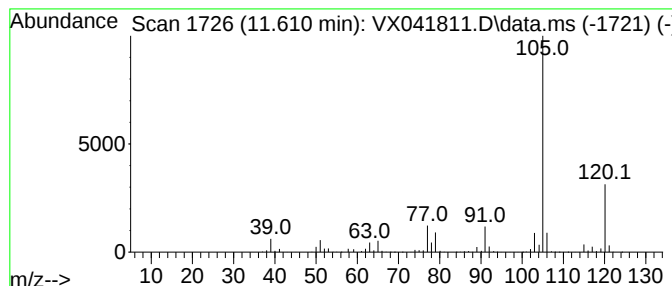
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-ethyl-4-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.610	112.90 ug/l	1680510	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			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061024\
Data File : VX041811.D
Acq On : 10 Jun 2024 15:31
Operator : JC/MD
Sample : P2767-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

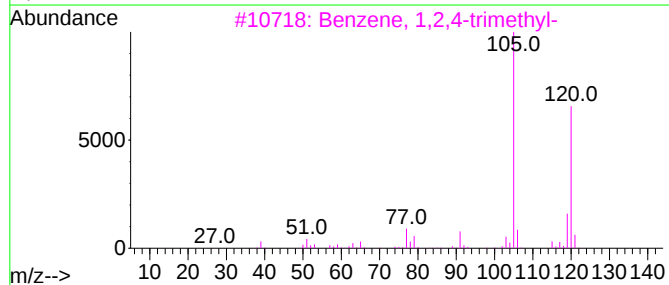
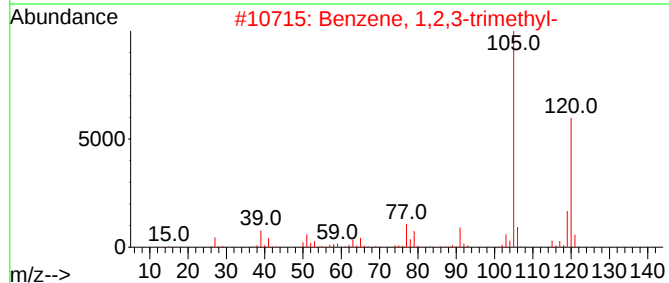
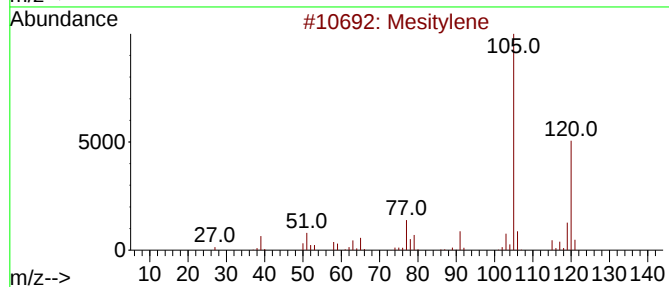
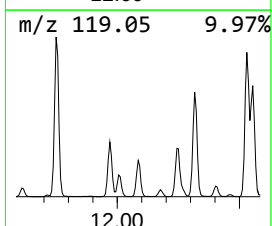
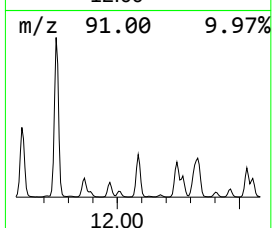
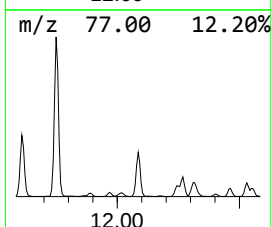
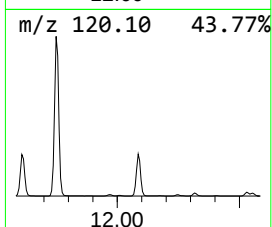
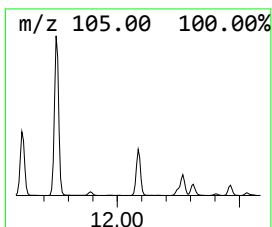
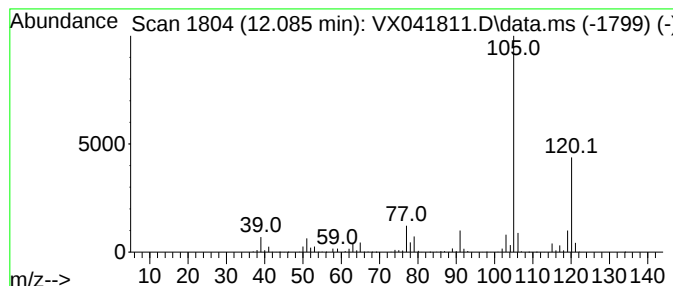
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethyl-3-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.085	85.09 ug/l	1266470	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	97
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061024\
Data File : VX041811.D
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Operator : JC/MD
Sample : P2767-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

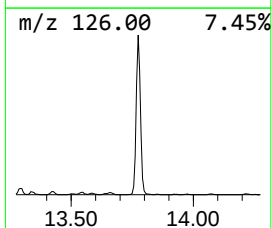
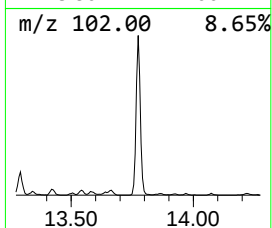
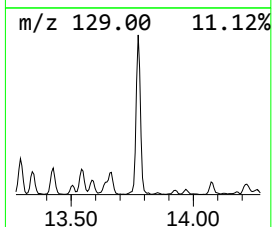
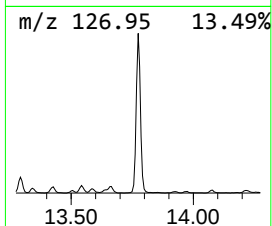
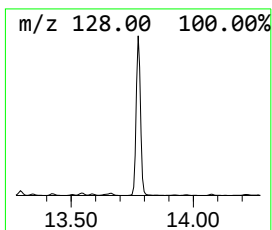
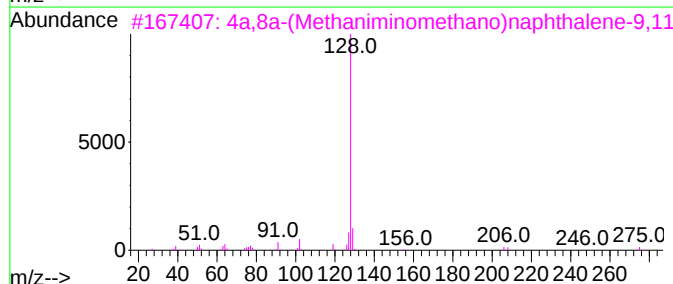
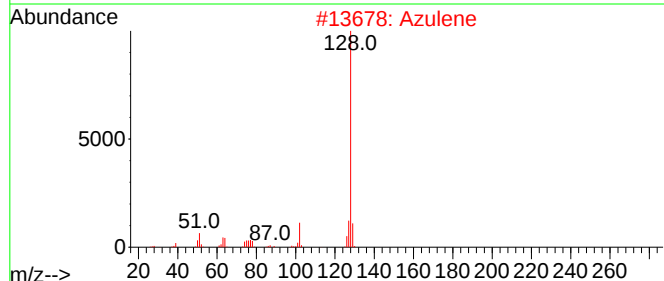
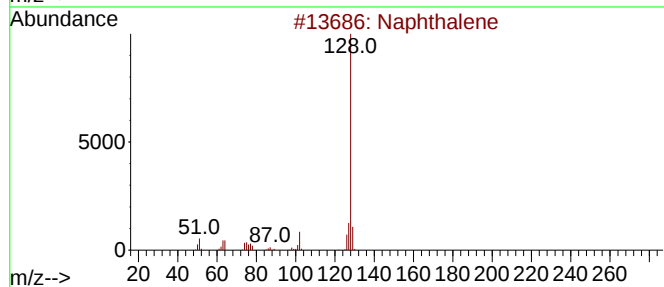
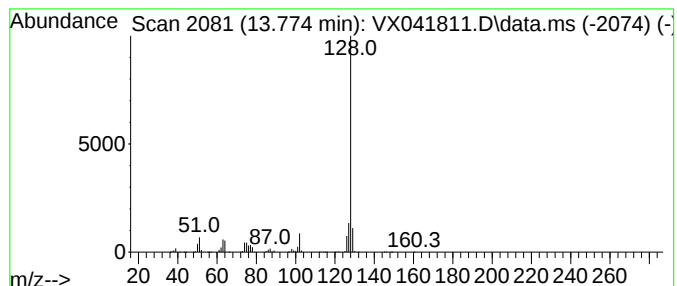
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Quant Title : SW846 8260

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

Peak Number 10 Azulene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.774	85.58 ug/l	1273860	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	95
2			Azulene	128	C10H8	000275-51-4	93
3			4a,8a-(Methaniminomethano)naphth...	275	C18H13NO2	069915-10-2	78
4			1H-Indene, 1-methylene-	128	C10H8	002471-84-3	47
5			1,3-Dicyanobenzene	128	C8H4N2	000626-17-5	45



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061024\
Data File : VX041811.D
Acq On : 10 Jun 2024 15:31
Operator : JC/MD
Sample : P2767-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X052924W.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-	1.721	189.7	ug/l	731651	1	5.550	192826	50.0
Pentane	1.935	122.5	ug/l	472357	1	5.550	192826	50.0
Pentane, 2-methyl-	2.813	158.3	ug/l	610664	1	5.550	192826	50.0
Pentane, 3-methyl-	3.093	161.1	ug/l	621440	1	5.550	192826	50.0
Cyclopentane, m...	4.294	394.4	ug/l	1520960	1	5.550	192826	50.0
Cyclopentane, 1...	6.153	99.2	ug/l	382638	1	5.550	192826	50.0
Benzene, 1-ethy...	11.378	212.5	ug/l	3162330	4	12.024	744220	50.0
Benzene, 1-ethy...	11.610	112.9	ug/l	1680510	4	12.024	744220	50.0
Benzene, 1-ethy...	12.085	85.1	ug/l	1266470	4	12.024	744220	50.0
Azulene	13.774	85.6	ug/l	1273860	4	12.024	744220	50.0