

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112525\
 Data File : VX048680.D
 Acq On : 25 Nov 2025 13:21
 Operator : JC/MD
 Sample : Q3716-01 20X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW5DL

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\82X110725W.M
 Title : SW846 8260

Signal : TIC: VX048680.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.276	27	31	43	rBV	121430	137144	9.85%	0.785%
2	1.380	43	48	55	rVV	377090	443909	31.88%	2.541%
3	1.435	55	57	65	rVB2	17572	22894	1.64%	0.131%
4	1.758	104	110	132	rBV	551143	879862	63.19%	5.037%
5	1.916	132	136	139	rBV2	10714	14795	1.06%	0.085%
6	1.965	139	144	157	rVV3	179915	346115	24.86%	1.981%
7	2.075	157	162	172	rVV	58962	93893	6.74%	0.538%
8	2.172	173	178	182	rBV	29808	45004	3.23%	0.258%
9	2.221	182	186	199	rVB	126213	218758	15.71%	1.252%
10	2.788	267	279	281	rBV4	35804	83754	6.02%	0.479%
11	2.831	281	286	289	rVV	71010	151100	10.85%	0.865%
12	2.873	289	293	315	rVB3	57095	168031	12.07%	0.962%
13	3.105	321	331	349	rBV	48814	125327	9.00%	0.717%
14	3.325	356	367	379	rBV2	13679	36460	2.62%	0.209%
15	3.446	379	387	398	rBV3	20204	50276	3.61%	0.288%
16	3.581	400	409	416	rBV3	9386	25844	1.86%	0.148%
17	3.660	416	422	427	rBV4	6759	15565	1.12%	0.089%
18	3.727	427	433	443	rVB	24778	57755	4.15%	0.331%
19	3.837	444	451	457	rBV3	14577	36682	2.63%	0.210%
20	4.099	485	494	503	rBV3	20855	60642	4.36%	0.347%
21	4.300	517	527	541	rBV2	71711	220499	15.84%	1.262%
22	4.416	541	546	558	rVB5	8253	23727	1.70%	0.136%
23	4.550	558	568	581	rVB3	3787	15556	1.12%	0.089%
24	5.184	662	672	682	rBV2	16796	54068	3.88%	0.310%
25	5.373	692	703	714	rBV2	121106	395183	28.38%	2.262%
26	5.538	722	730	763	rVB	186230	638316	45.84%	3.654%
27	5.812	766	775	786	rBV5	11770	37331	2.68%	0.214%
28	5.946	786	797	805	rVV	123621	348613	25.04%	1.996%
29	6.025	807	810	822	rVV	24037	65857	4.73%	0.377%
30	6.153	823	831	835	rVV7	9566	31918	2.29%	0.183%
31	6.239	836	845	855	rVV2	29056	105570	7.58%	0.604%
32	6.342	856	862	874	rVB5	7408	23612	1.70%	0.135%
33	6.598	894	904	912	rBV6	6479	20479	1.47%	0.117%
34	6.745	919	928	939	rBV	349400	901213	64.73%	5.159%
35	7.153	989	995	1005	rBV3	8084	19783	1.42%	0.113%
36	7.360	1019	1029	1041	rBV6	24521	70832	5.09%	0.406%

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Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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37	7.970	1119	1129	1140	rVB2	20776	46214	3.32%	0.265%
38	8.092	1141	1149	1153	rBV	28459	60482	4.34%	0.346%
39	8.171	1159	1162	1171	rVB5	7839	14915	1.07%	0.085%
40	8.421	1197	1203	1210	rBV4	8319	15987	1.15%	0.092%
41	8.494	1210	1215	1222	rVB5	9368	17922	1.29%	0.103%
42	8.635	1222	1238	1245	rBV	708953	1182211	84.91%	6.768%
43	8.702	1245	1249	1257	rVB	67501	117574	8.44%	0.673%
44	10.037	1463	1468	1486	rVB	893279	1256836	90.27%	7.195%
45	10.177	1486	1491	1498	rBV	235703	315577	22.67%	1.807%
46	10.281	1502	1508	1517	rBV	934992	1392348	100.00%	7.971%
47	10.622	1559	1564	1575	rBV	335798	473574	34.01%	2.711%
48	10.945	1611	1617	1623	rBV	20463	26883	1.93%	0.154%
49	11.061	1631	1636	1648	rBV	685061	945317	67.89%	5.412%
50	11.287	1668	1673	1680	rBV	52124	68558	4.92%	0.392%
51	11.360	1680	1685	1693	rVV	379374	744796	53.49%	4.264%
52	11.433	1693	1697	1709	rVB	275309	352251	25.30%	2.017%
53	11.597	1716	1724	1732	rBV	216209	290068	20.83%	1.661%
54	11.732	1741	1746	1754	rBV	868169	1114867	80.07%	6.383%
55	11.951	1779	1782	1785	rBV	13587	15283	1.10%	0.087%
56	12.000	1785	1790	1797	rVV	929488	1250673	89.82%	7.160%
57	12.067	1797	1801	1810	rVB	209902	272589	19.58%	1.561%
58	12.225	1821	1827	1830	rBV2	125131	186589	13.40%	1.068%
59	12.299	1835	1839	1848	rVB2	99564	141987	10.20%	0.813%
60	12.390	1848	1854	1859	rBV7	8120	17361	1.25%	0.099%
61	12.445	1859	1863	1870	rVB	22998	29152	2.09%	0.167%
62	12.512	1870	1874	1876	rBV	55023	69247	4.97%	0.396%
63	12.536	1876	1878	1883	rVV	49674	54113	3.89%	0.310%
64	12.591	1883	1887	1891	rVV	87424	114562	8.23%	0.656%
65	12.628	1891	1893	1897	rVV	18518	19089	1.37%	0.109%
66	12.677	1897	1901	1911	rVV2	52651	81801	5.88%	0.468%
67	12.829	1922	1926	1931	rBV	22661	28558	2.05%	0.163%
68	12.902	1931	1938	1941	rBV	54208	70063	5.03%	0.401%
69	12.945	1941	1945	1951	rVB	85402	104607	7.51%	0.599%
70	13.024	1951	1958	1961	rBV3	9448	21115	1.52%	0.121%
71	13.146	1974	1978	1983	rVB	48968	61667	4.43%	0.353%
72	13.274	1989	1999	2004	rBV2	107921	167319	12.02%	0.958%
73	13.402	2016	2020	2028	rVB3	12219	19342	1.39%	0.111%
74	13.524	2037	2040	2043	rVV	15503	19879	1.43%	0.114%
75	13.561	2043	2046	2052	rVV3	18424	35542	2.55%	0.203%
76	13.646	2052	2060	2066	rVB2	13586	30460	2.19%	0.174%

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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

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Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X110725W.M
Title : SW846 8260

77	13.756	2073	2078	2085	rVB	71160	90950	6.53%	0.521%
78	14.054	2124	2127	2133	rVB2	12232	15493	1.11%	0.089%
79	14.195	2145	2150	2156	rBV3	11080	17128	1.23%	0.098%
80	14.615	2208	2219	2229	rVB	51453	69742	5.01%	0.399%
81	14.755	2237	2242	2255	rVB	30167	40311	2.90%	0.231%

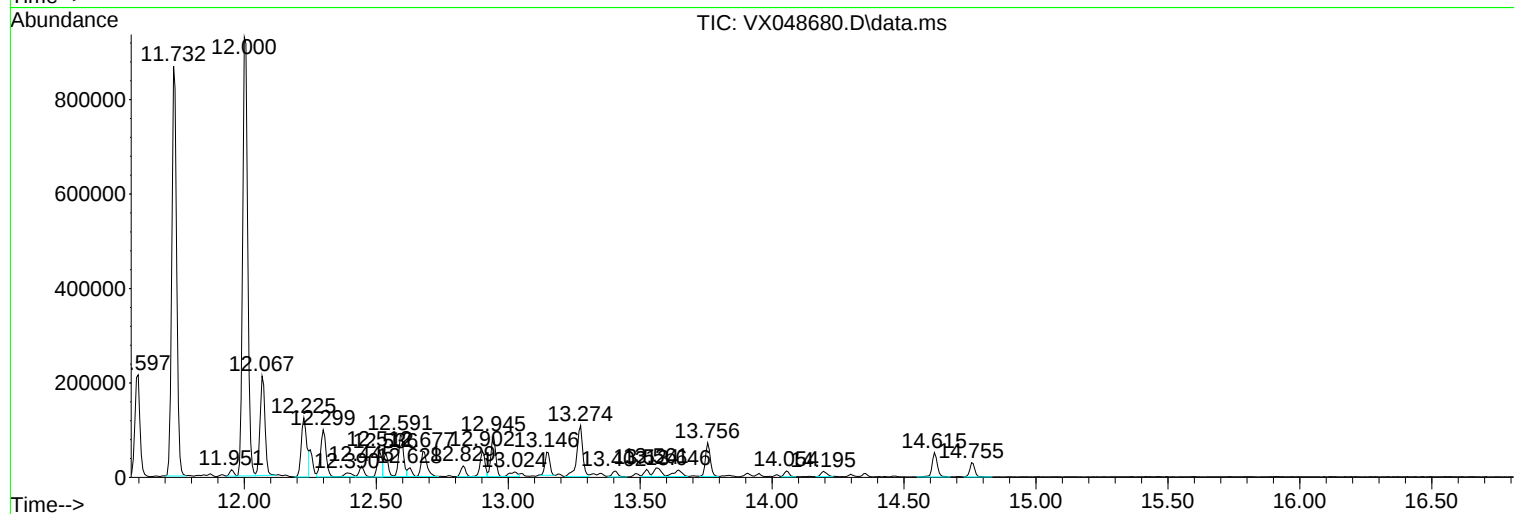
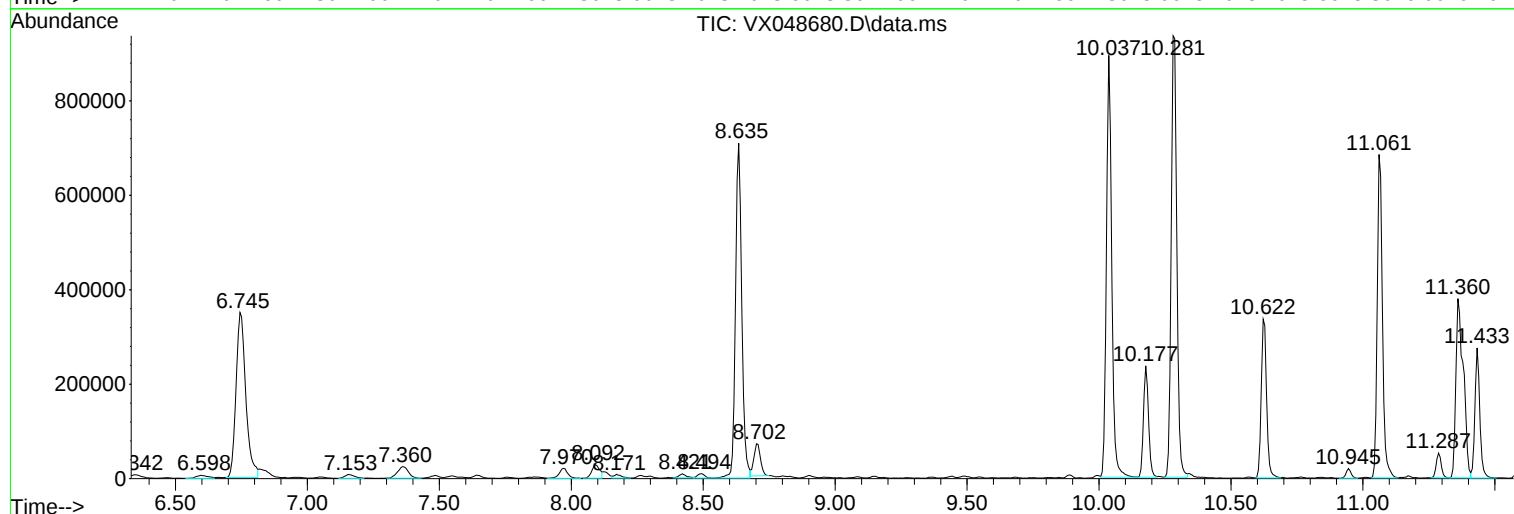
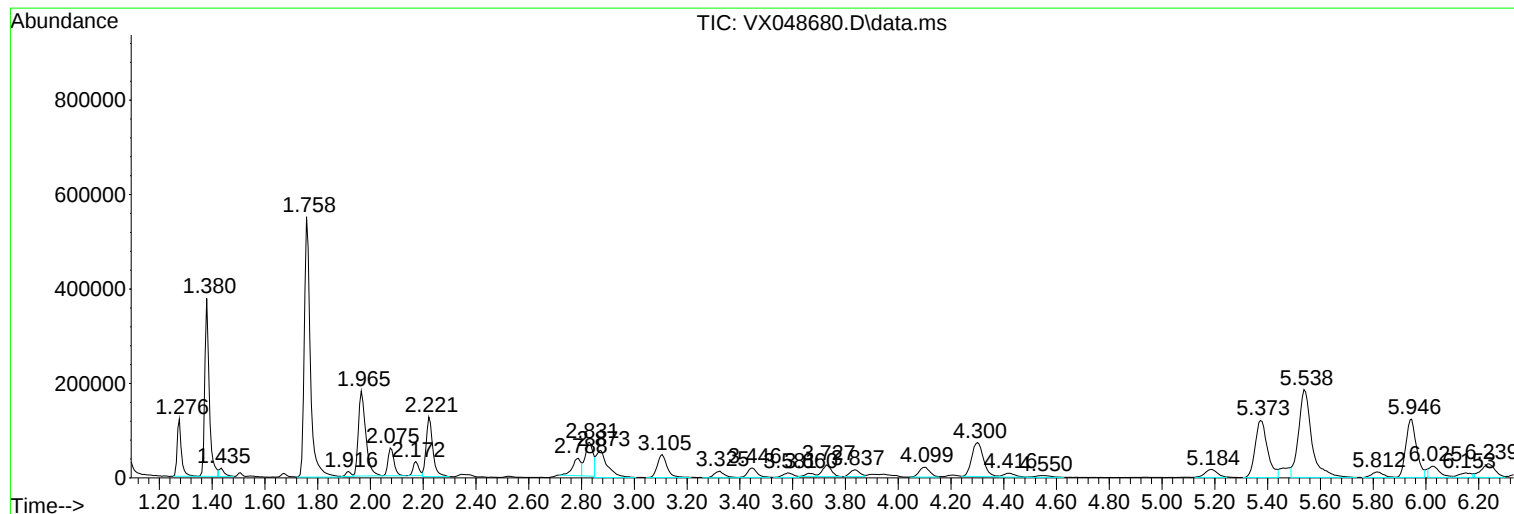
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Instrument :
MSVOA_X
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X110725W.M
Quant Title : SW846 8260

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



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Instrument :
MSVOA_X
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MW5DL

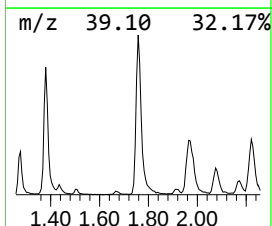
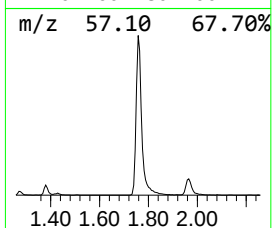
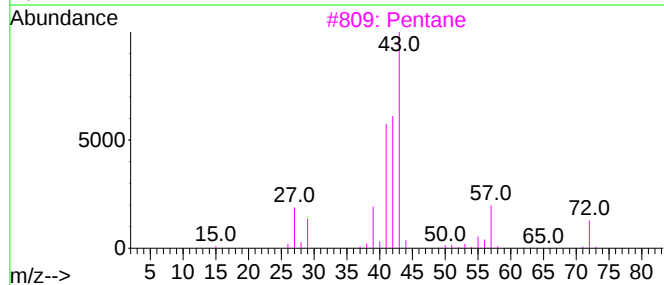
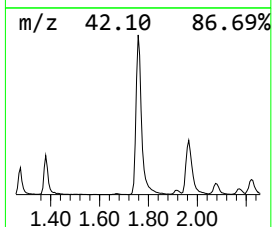
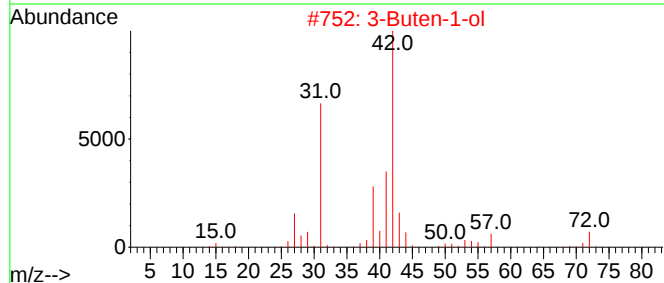
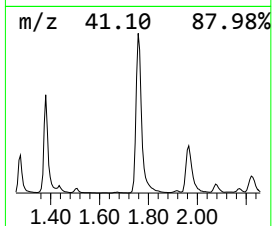
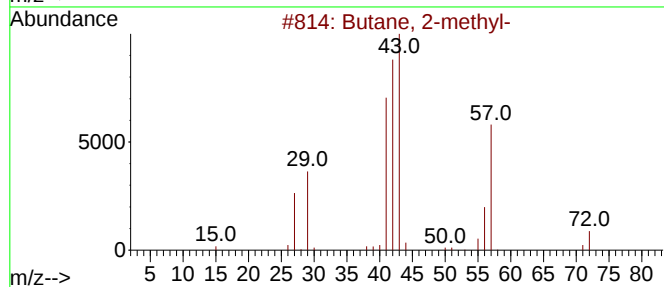
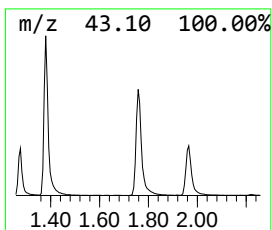
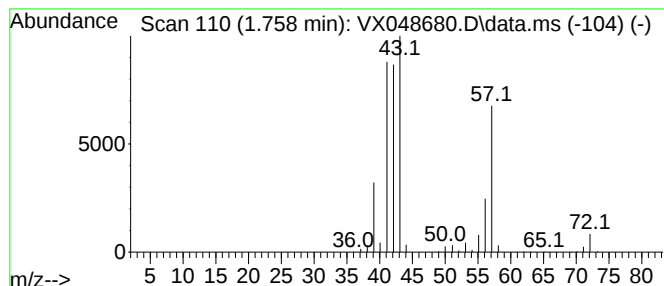
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X110725W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.758	68.92 ug/l	879862	Pentafluorobenzene	5.538

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methyl-	72	C5H12	000078-78-4	91
2			3-Buten-1-ol	72	C4H8O	000627-27-0	47
3			Pentane	72	C5H12	000109-66-0	45
4			1-Butene	56	C4H8	000106-98-9	14
5			Azetidine	57	C3H7N	000503-29-7	9



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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW5DL

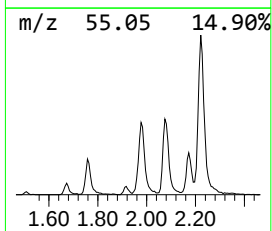
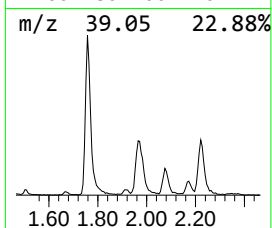
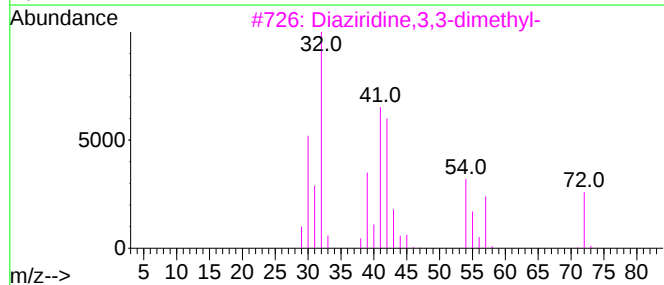
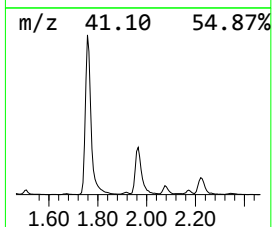
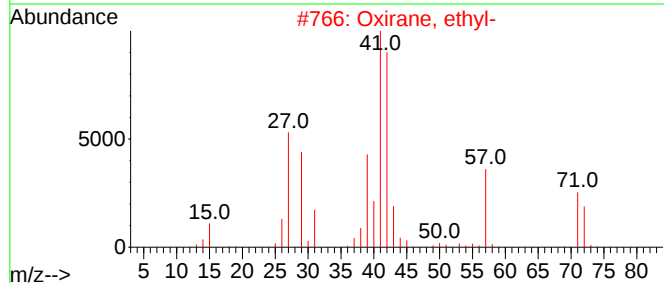
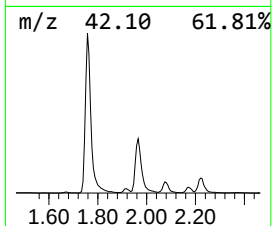
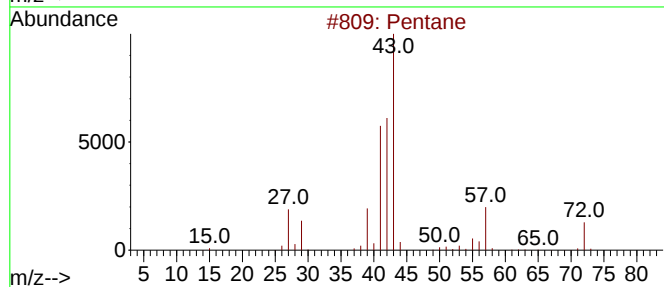
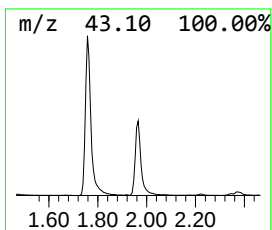
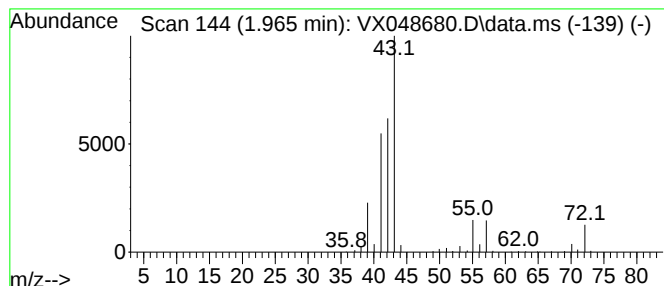
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X110725W.M
Quant Title : SW846 8260

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

Peak Number 4 Pentane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.965	27.11 ug/l	346115	Pentafluorobenzene	5.538

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane	72	C5H12	000109-66-0	87
2			Oxirane, ethyl-	72	C4H8O	000106-88-7	43
3			Diaziridine,3,3-dimethyl-	72	C3H8N2	004901-76-2	35
4			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	17
5			Butane, 2-methyl-	72	C5H12	000078-78-4	9



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Instrument :
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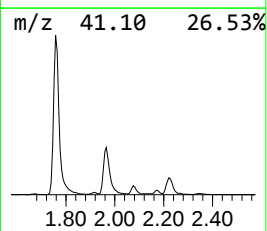
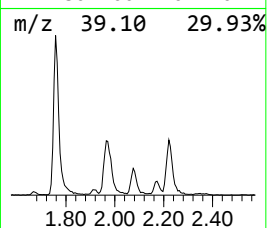
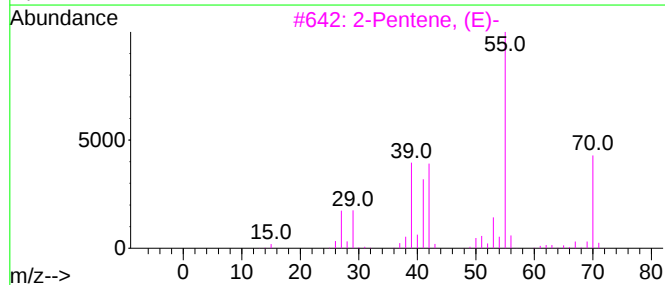
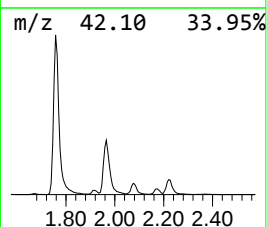
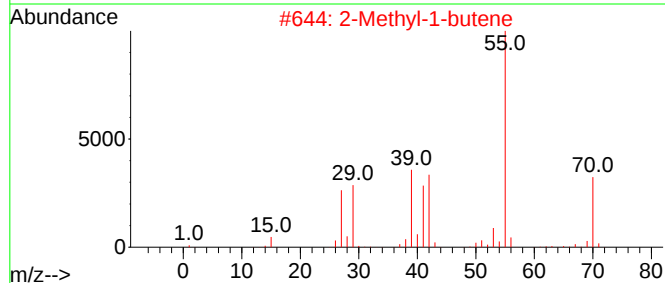
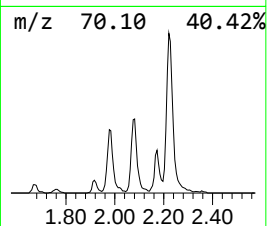
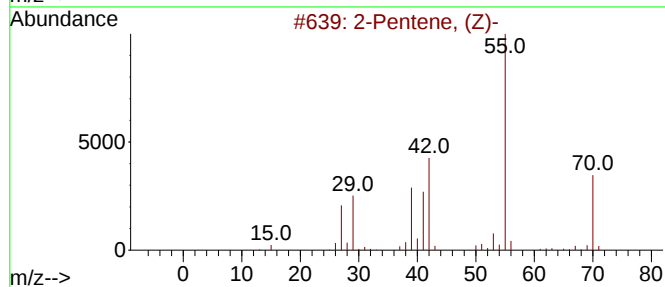
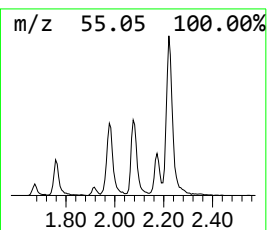
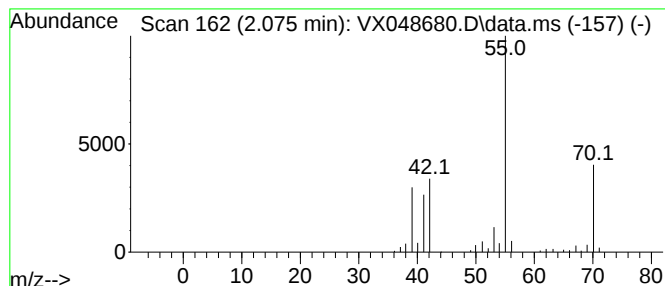
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X110725W.M
Quant Title : SW846 8260

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

Peak Number 5 2-Pentene, (Z)- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.075	7.35 ug/l	93893	Pentafluorobenzene	5.538

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentene, (Z)-	70	C5H10	000627-20-3	91
2			2-Methyl-1-butene	70	C5H10	000563-46-2	91
3			2-Pentene, (E)-	70	C5H10	000646-04-8	90
4			Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	90
5			Cyclopropane, 1,1-dimethyl-	70	C5H10	001630-94-0	87



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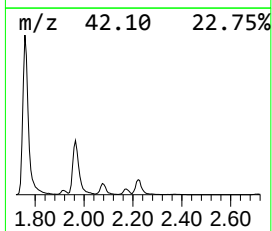
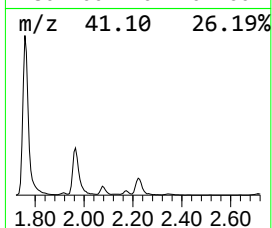
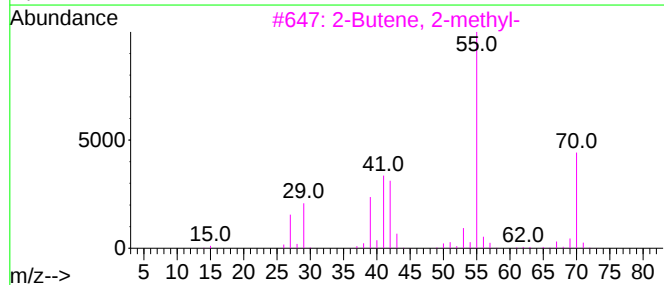
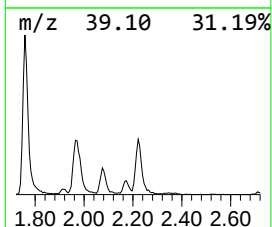
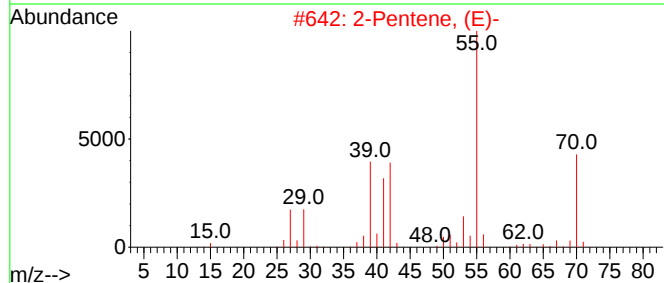
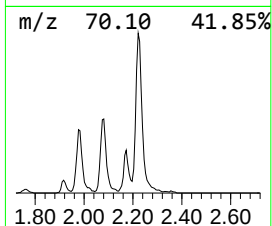
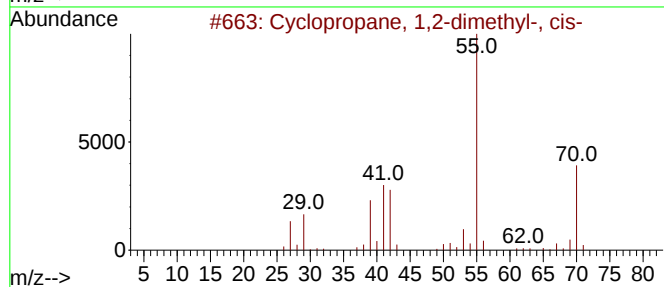
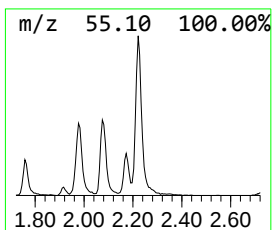
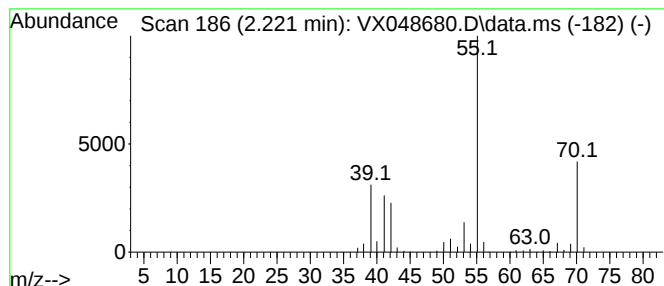
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Quant Title : SW846 8260

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

Peak Number 6 Cyclopropane, 1,2-dimethyl-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.221	17.14 ug/l	218758	Pentafluorobenzene	5.538

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	90
2			2-Pentene, (E)-	70	C5H10	000646-04-8	87
3			2-Butene, 2-methyl-	70	C5H10	000513-35-9	86
4			Cyclopropane, 1,1-dimethyl-	70	C5H10	001630-94-0	80
5			Cyclopropane, 1,2-dimethyl-, trans-	70	C5H10	002402-06-4	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112525\
Data File : VX048680.D
Acq On : 25 Nov 2025 13:21
Operator : JC/MD
Sample : Q3716-01 20X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW5DL

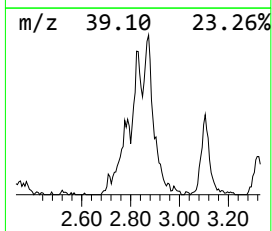
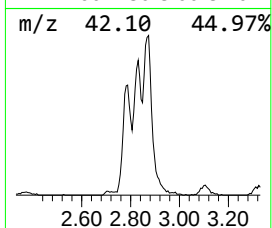
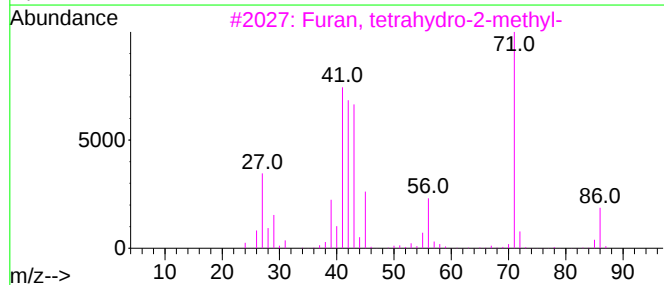
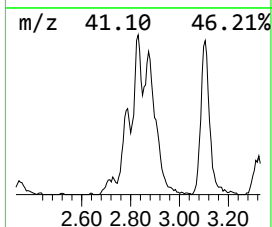
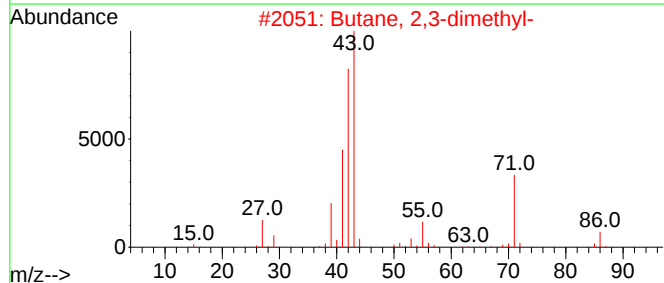
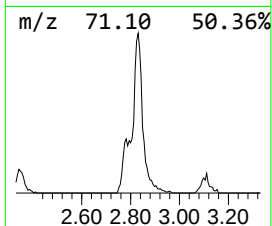
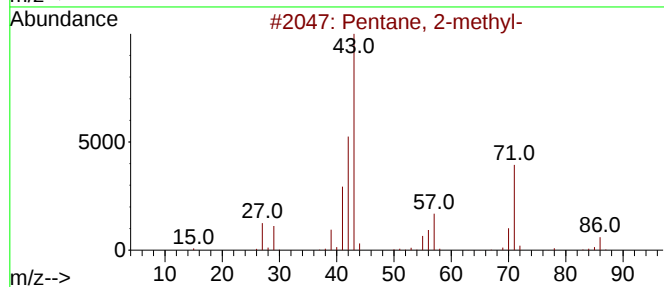
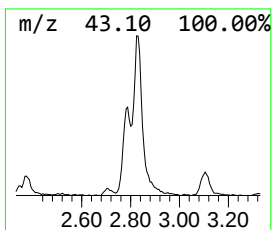
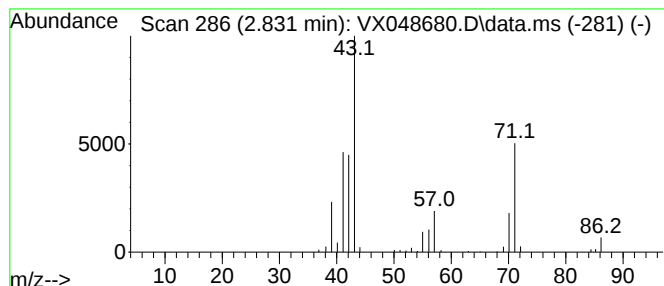
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Quant Title : SW846 8260

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

Peak Number 7 Pentane, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.831	11.84 ug/l	151100	Pentafluorobenzene	5.538

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			Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	38
4			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	38
5			Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	37



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112525\
Data File : VX048680.D
Acq On : 25 Nov 2025 13:21
Operator : JC/MD
Sample : Q3716-01 20X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW5DL

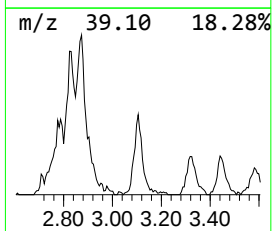
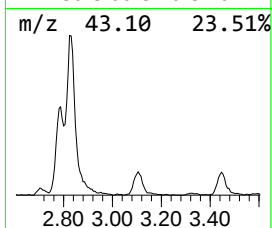
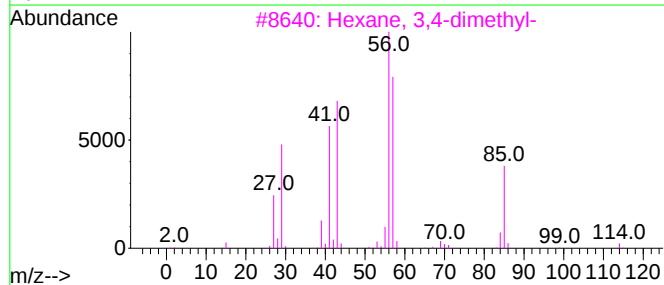
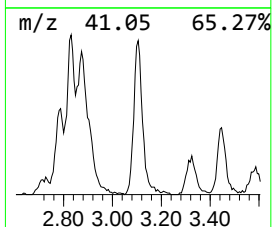
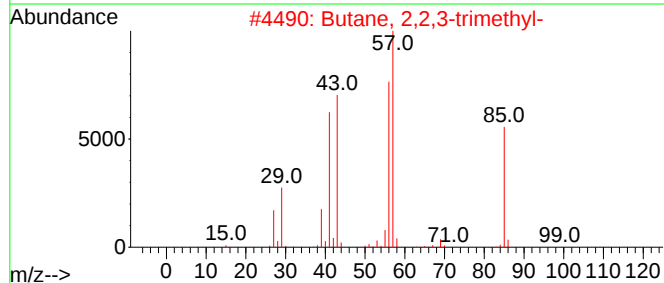
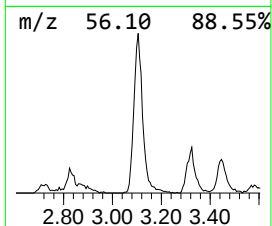
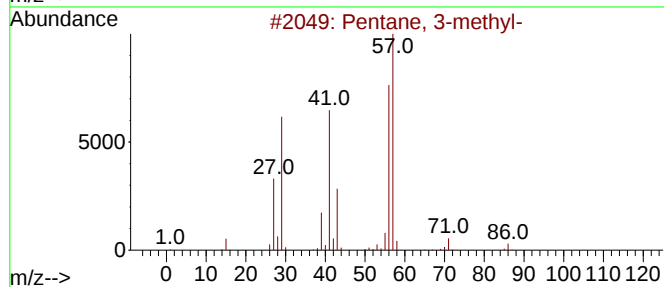
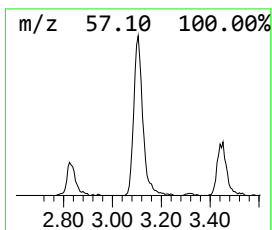
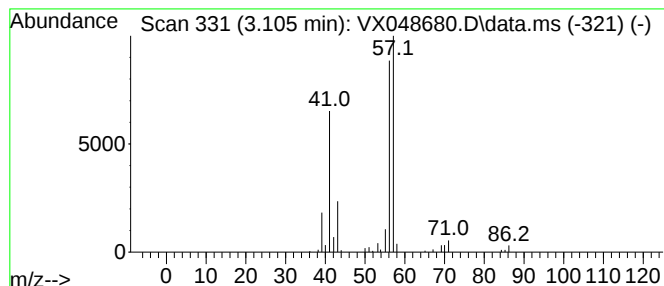
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Quant Title : SW846 8260

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

Peak Number 9 Pentane, 3-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.105	9.82 ug/l	125327	Pentafluorobenzene	5.538

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-methyl-	86	C6H14	000096-14-0	90
2			Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	83
3			Hexane, 3,4-dimethyl-	114	C8H18	000583-48-2	78
4			Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	78
5			Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112525\
Data File : VX048680.D
Acq On : 25 Nov 2025 13:21
Operator : JC/MD
Sample : Q3716-01 20X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW5DL

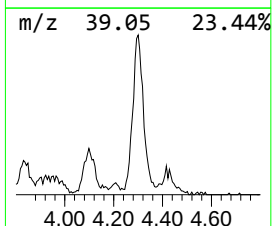
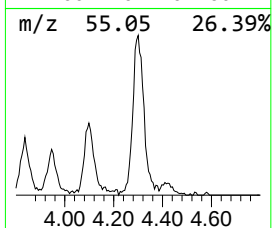
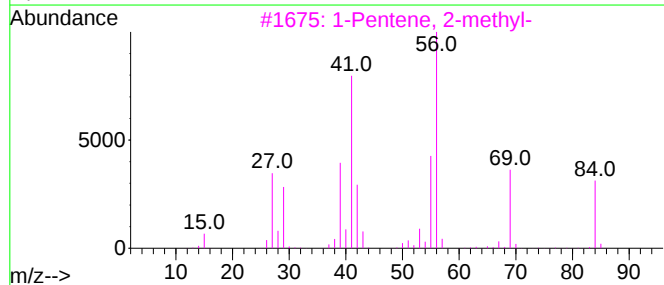
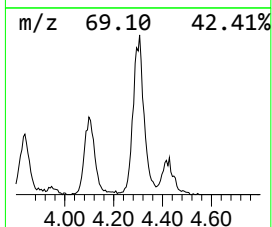
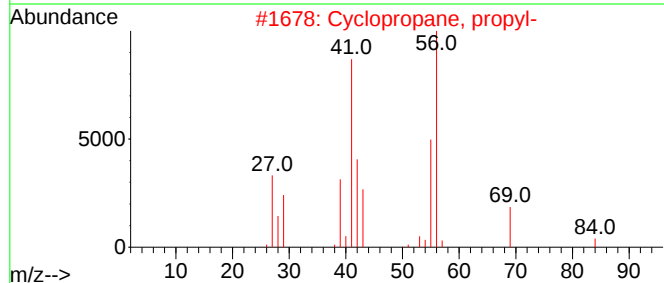
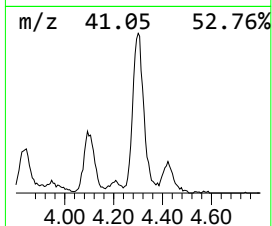
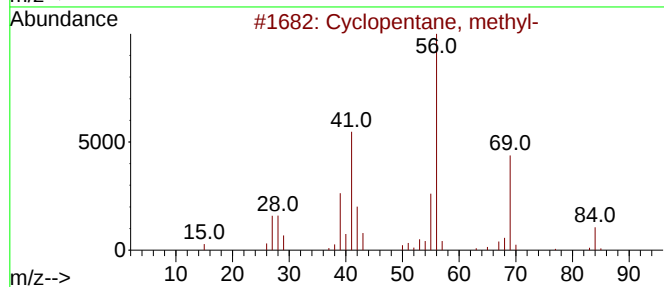
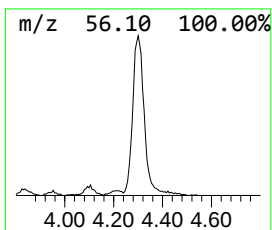
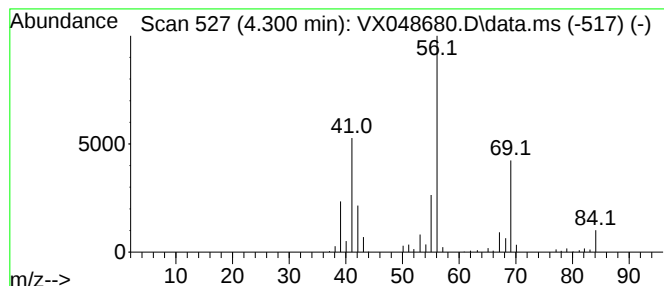
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Quant Title : SW846 8260

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

Peak Number 10 Cyclopentane, methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.300	17.27 ug/l	220499	Pentafluorobenzene	5.538

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2		Cyclopropane, propyl-	84	C6H12	002415-72-7	64
3		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	59
4		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	50
5		1-Pentanol, 3-methyl-	102	C6H14O	000589-35-5	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112525\
Data File : VX048680.D
Acq On : 25 Nov 2025 13:21
Operator : JC/MD
Sample : Q3716-01 20X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW5DL

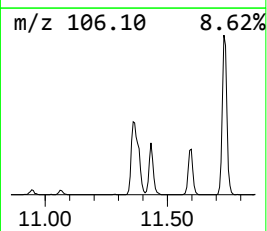
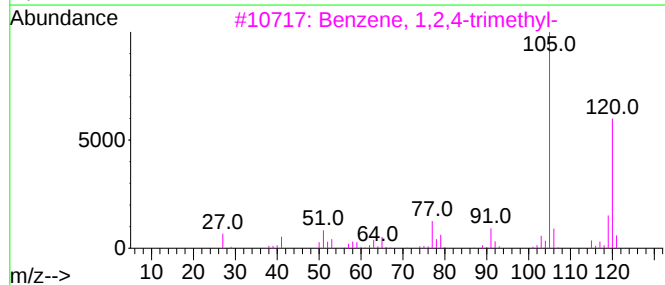
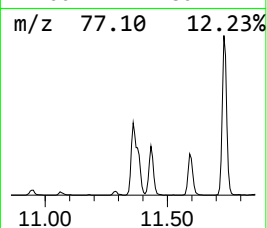
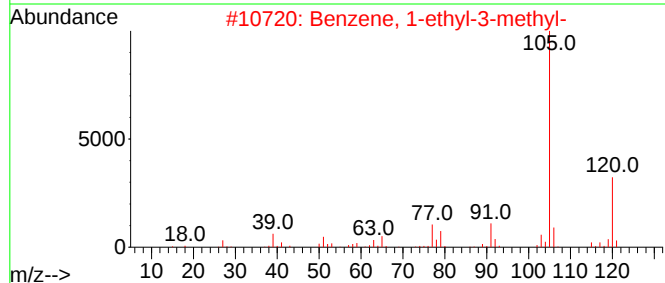
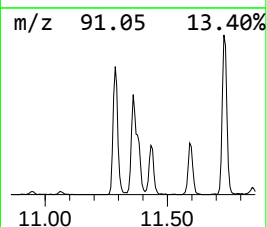
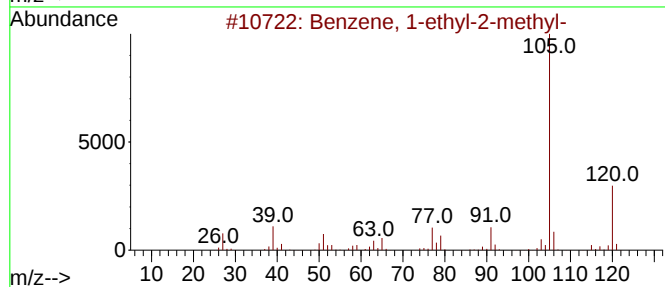
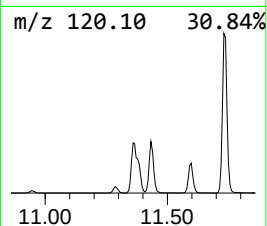
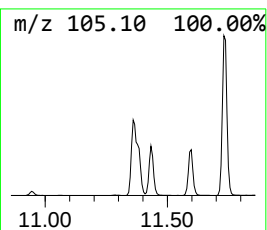
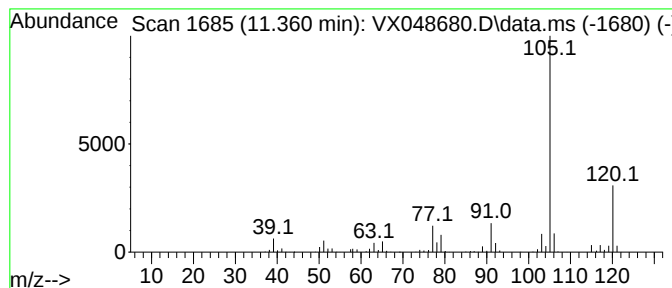
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Quant Title : SW846 8260

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

Peak Number 11 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.360	29.78 ug/l	744796	1,4-Dichlorobenzene-d4	12.006

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	94
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112525\
Data File : VX048680.D
Acq On : 25 Nov 2025 13:21
Operator : JC/MD
Sample : Q3716-01 20X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW5DL

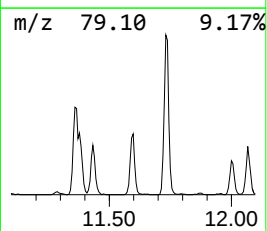
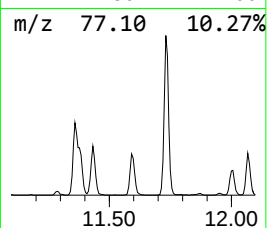
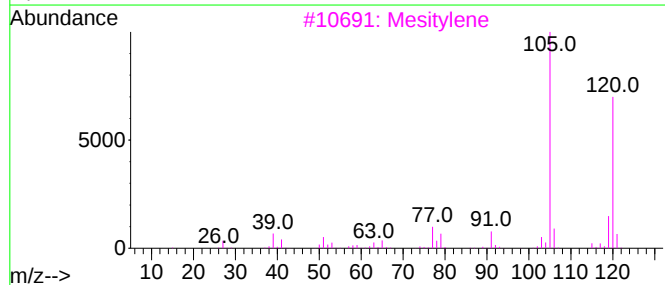
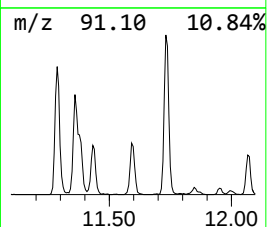
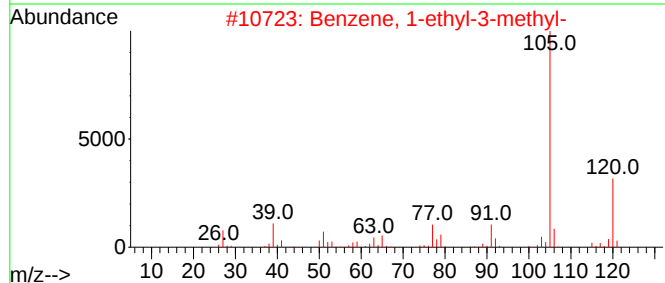
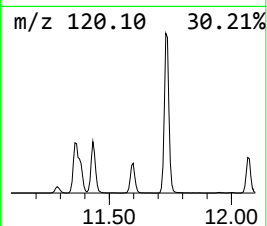
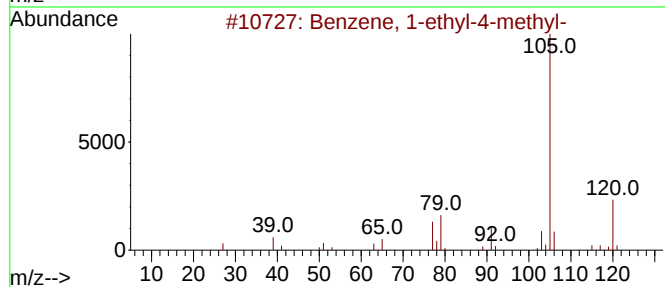
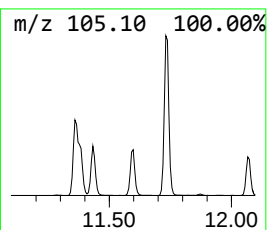
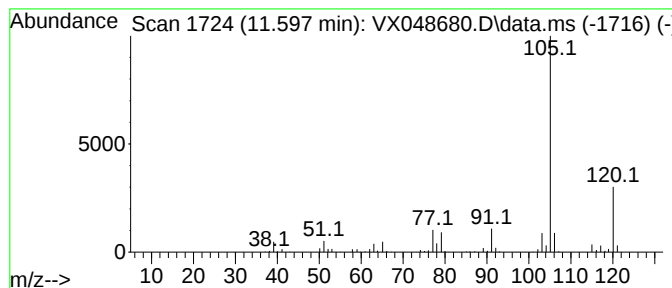
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Quant Title : SW846 8260

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

Peak Number 12 Benzene, 1-ethyl-4-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.597	11.60 ug/l	290068	1,4-Dichlorobenzene-d4	12.006

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112525\
Data File : VX048680.D
Acq On : 25 Nov 2025 13:21
Operator : JC/MD
Sample : Q3716-01 20X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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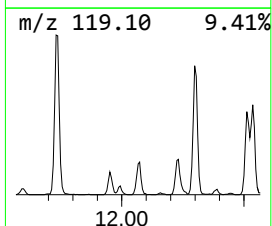
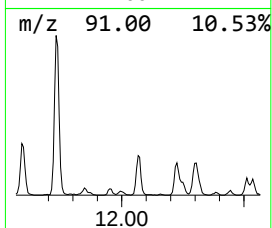
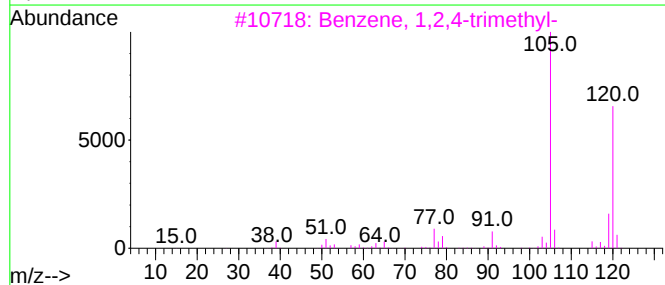
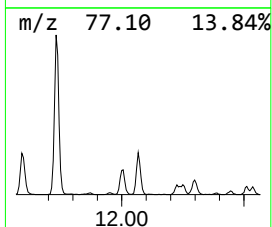
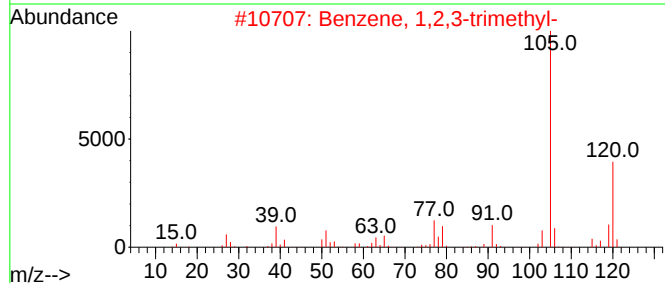
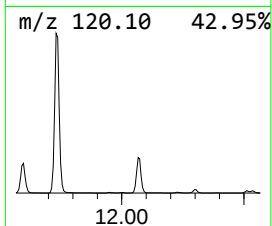
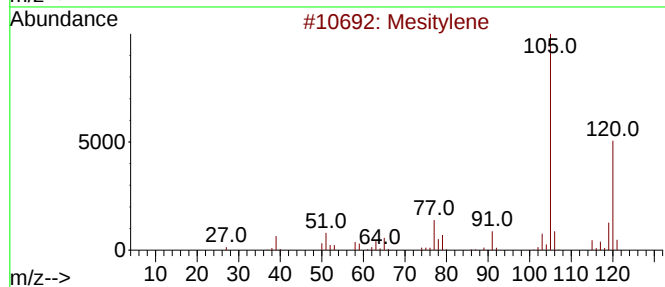
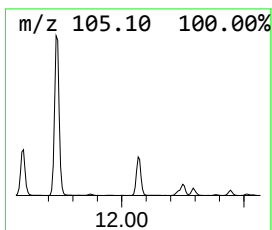
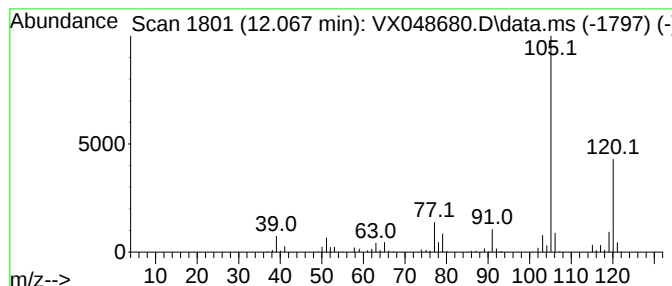
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Quant Title : SW846 8260

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

Peak Number 13 Benzene, 1,2,3-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.067	10.90 ug/l	272589	1,4-Dichlorobenzene-d4	12.006

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	94
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112525\
Data File : VX048680.D
Acq On : 25 Nov 2025 13:21
Operator : JC/MD
Sample : Q3716-01 20X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

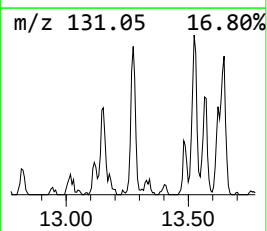
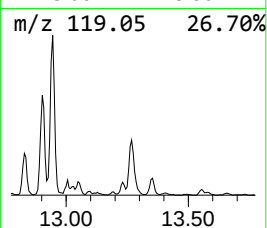
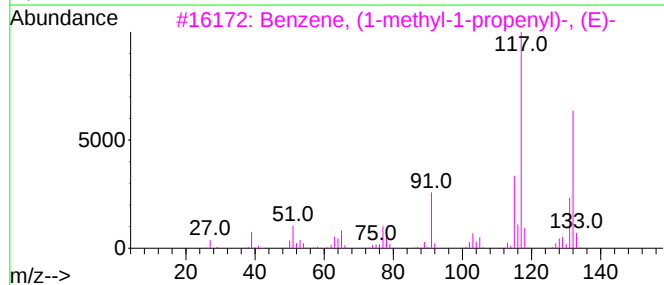
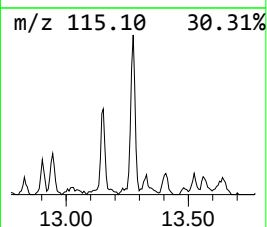
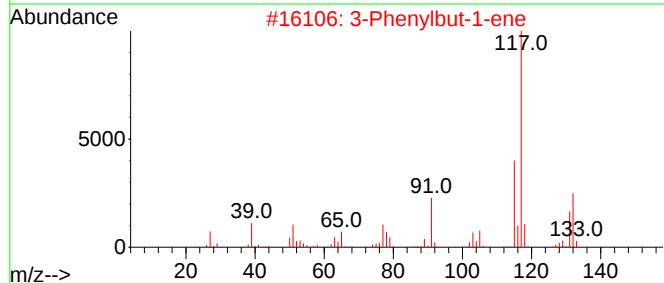
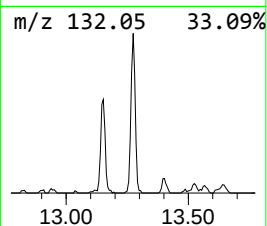
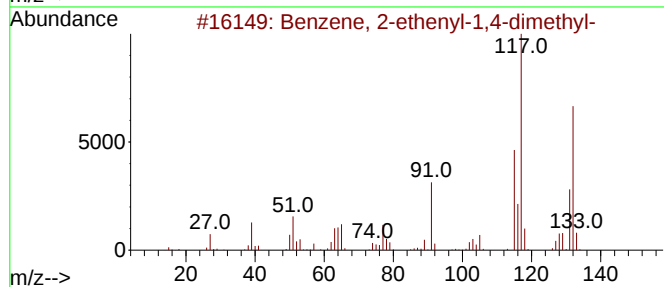
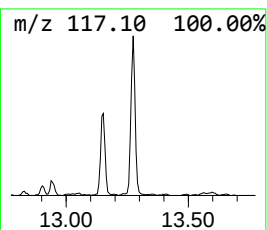
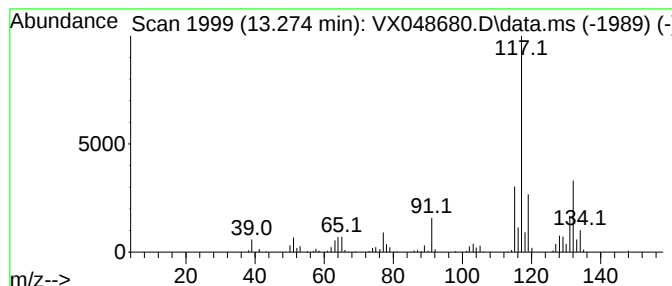
Instrument :
MSVOA_X
ClientSampleId :
MW5DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X110725W.M
Quant Title : SW846 8260

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

Peak Number 15 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD			R.T.
13.274	6.69 ug/l	167319	1,4-Dichlorobenzene-d4			12.006
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-		132	C10H12	002039-89-6	90
2	3-Phenylbut-1-ene		132	C10H12	000934-10-1	81
3	Benzene, (1-methyl-1-propenyl)-,...		132	C10H12	000768-00-3	81
4	Benzene, 1-ethenyl-3-ethyl-		132	C10H12	007525-62-4	76
5	1H-Indene, 2,3-dihydro-4-methyl-		132	C10H12	000824-22-6	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112525\
Data File : VX048680.D
Acq On : 25 Nov 2025 13:21
Operator : JC/MD
Sample : Q3716-01 20X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW5DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X110725W.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-	1.758	68.9	ug/l	879862	1	5.538	638316	50.0
Pentane	1.965	27.1	ug/l	346115	1	5.538	638316	50.0
2-Pentene, (Z)-	2.075	7.3	ug/l	93893	1	5.538	638316	50.0
Cyclopropane, 1...	2.221	17.1	ug/l	218758	1	5.538	638316	50.0
Pentane, 2-methyl-	2.831	11.8	ug/l	151100	1	5.538	638316	50.0
Pentane, 3-methyl-	3.105	9.8	ug/l	125327	1	5.538	638316	50.0
Cyclopentane, m...	4.300	17.3	ug/l	220499	1	5.538	638316	50.0
Benzene, 1-ethy...	11.360	29.8	ug/l	744796	4	12.006	1250670	50.0
Benzene, 1-ethy...	11.597	11.6	ug/l	290068	4	12.006	1250670	50.0
Benzene, 1,2,3-...	12.067	10.9	ug/l	272589	4	12.006	1250670	50.0
Benzene, 2-ethe...	13.274	6.7	ug/l	167319	4	12.006	1250670	50.0