

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061522\
 Data File : VX029507.D
 Acq On : 15 Jun 2022 14:38
 Operator : JC/MD
 Sample : N3160-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-2

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\82X060722W.M
 Title : SW846 8260

Signal : TIC: VX029507.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.142	6	9	15	rVB	49639	42730	2.23%	0.251%
2	2.337	200	205	209	rBV2	10820	19191	1.00%	0.113%
3	2.782	268	278	292	rBV	48904	119392	6.24%	0.701%
4	3.099	320	330	338	rBV2	11122	25041	1.31%	0.147%
5	4.056	475	487	496	rBV2	7661	22242	1.16%	0.131%
6	4.214	500	513	524	rBV	91886	275231	14.38%	1.615%
7	4.458	543	553	564	rVB4	6625	23062	1.20%	0.135%
8	5.111	649	660	674	rBV7	15710	66452	3.47%	0.390%
9	5.391	691	706	719	rBV	209451	612161	31.98%	3.593%
10	5.556	719	733	740	rVV	271241	808312	42.22%	4.744%
11	5.623	740	744	758	rVB2	118523	320618	16.75%	1.882%
12	5.958	788	799	812	rBV	225139	608123	31.77%	3.569%
13	6.251	825	847	860	rBV	553889	1689354	88.24%	9.915%
14	6.763	920	931	942	rBV	397921	953007	49.78%	5.593%
15	6.964	958	964	976	rBV2	9319	23482	1.23%	0.138%
16	7.128	981	991	1004	rVB4	15102	38858	2.03%	0.228%
17	7.366	1021	1030	1041	rVB4	16411	39967	2.09%	0.235%
18	7.500	1044	1052	1057	rBV2	32032	69918	3.65%	0.410%
19	7.561	1057	1062	1066	rVV2	45711	98139	5.13%	0.576%
20	7.604	1066	1069	1077	rVB2	33662	63895	3.34%	0.375%
21	7.775	1090	1097	1106	rVB2	27163	57555	3.01%	0.338%
22	7.988	1123	1132	1143	rVB	451527	882625	46.10%	5.180%
23	8.110	1143	1152	1160	rBV	454007	886401	46.30%	5.203%
24	8.317	1179	1186	1192	rVB5	10815	21662	1.13%	0.127%
25	8.390	1192	1198	1202	rBV2	17373	31885	1.67%	0.187%
26	8.653	1225	1241	1249	rBV	1155524	1914415	100.00%	11.236%
27	8.982	1291	1295	1300	rVB3	18341	28624	1.50%	0.168%
28	9.104	1307	1315	1320	rVB3	11575	23991	1.25%	0.141%
29	9.165	1320	1325	1328	rVV	19286	29739	1.55%	0.175%
30	9.201	1328	1331	1336	rVB2	18546	26760	1.40%	0.157%
31	9.512	1376	1382	1388	rBV5	13017	30014	1.57%	0.176%
32	9.835	1430	1435	1439	rBV2	12982	19556	1.02%	0.115%
33	10.055	1466	1471	1477	rBV	793608	1041696	54.41%	6.114%
34	10.280	1503	1508	1510	rVV3	16722	28586	1.49%	0.168%
35	10.305	1510	1512	1518	rVB4	17220	22657	1.18%	0.133%
36	10.591	1548	1559	1566	rBV7	17916	47209	2.47%	0.277%

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Integrator: RTE

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37	10.780	1581	1590	1594	rBV4	20334	40174	2.10%	0.236%
38	11.085	1634	1640	1644	rBV	900603	1152277	60.19%	6.763%
39	11.122	1644	1646	1651	rVV2	26470	37896	1.98%	0.222%
40	11.219	1655	1662	1667	rVB2	29919	43923	2.29%	0.258%

41	11.622	1720	1728	1734	rBV6	11809	26186	1.37%	0.154%
42	11.865	1761	1768	1770	rBV	50569	74867	3.91%	0.439%
43	11.890	1770	1772	1776	rVV	47391	58735	3.07%	0.345%
44	12.024	1789	1794	1805	rBV	767662	971154	50.73%	5.700%
45	12.122	1805	1810	1813	rBV4	12415	20979	1.10%	0.123%

46	12.250	1827	1831	1835	rVB3	17024	20539	1.07%	0.121%
47	12.317	1835	1842	1849	rBV2	65036	128357	6.70%	0.753%
48	12.408	1853	1857	1863	rBV	32356	46664	2.44%	0.274%
49	12.475	1863	1868	1872	rVB	40666	53800	2.81%	0.316%
50	12.603	1882	1889	1893	rBV4	15999	31954	1.67%	0.188%

51	12.725	1902	1909	1913	rBV4	27769	56787	2.97%	0.333%
52	12.762	1913	1915	1920	rVB2	17965	21830	1.14%	0.128%
53	12.835	1920	1927	1933	rVB3	46652	95929	5.01%	0.563%
54	12.902	1933	1938	1939	rBV2	55049	66878	3.49%	0.393%
55	12.920	1939	1941	1947	rVB2	69692	93296	4.87%	0.548%

56	12.987	1947	1952	1955	rBV4	18973	23622	1.23%	0.139%
57	13.030	1955	1959	1966	rVB3	35855	62511	3.27%	0.367%
58	13.115	1966	1973	1974	rBV2	40145	61029	3.19%	0.358%
59	13.140	1974	1977	1982	rVB	86665	107595	5.62%	0.632%
60	13.194	1982	1986	1991	rBV	32518	46088	2.41%	0.271%

61	13.249	1991	1995	1998	rBV3	14501	23776	1.24%	0.140%
62	13.292	1998	2002	2008	rVB4	35953	62197	3.25%	0.365%
63	13.420	2022	2023	2034	rVB5	22819	36422	1.90%	0.214%
64	13.518	2034	2039	2041	rBV3	31053	50624	2.64%	0.297%
65	13.579	2046	2049	2053	rBV3	85792	119495	6.24%	0.701%

66	13.621	2053	2056	2061	rVV2	49100	87295	4.56%	0.512%
67	13.664	2061	2063	2068	rVB2	42489	62997	3.29%	0.370%
68	13.755	2072	2078	2081	rBV	50259	70346	3.67%	0.413%
69	13.798	2081	2085	2090	rVB3	44629	64789	3.38%	0.380%
70	13.853	2090	2094	2099	rVB2	80282	113974	5.95%	0.669%

71	13.926	2099	2106	2110	rBV4	95607	178323	9.31%	1.047%
72	13.975	2110	2114	2118	rBV2	28917	41760	2.18%	0.245%
73	14.072	2126	2130	2133	rBV3	38046	54872	2.87%	0.322%
74	14.139	2138	2141	2143	rVV3	18090	23881	1.25%	0.140%
75	14.164	2143	2145	2149	rVV4	18315	28410	1.48%	0.167%

76	14.213	2149	2153	2164	rVB2	96782	159619	8.34%	0.937%
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Instrument :
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ClientSampleId :
MW-2

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

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Min Area: 3 % of largest Peak

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77	14.322	2164	2171	2176	rBV	77599	121176	6.33%	0.711%
78	14.371	2176	2179	2184	rVB2	68970	91107	4.76%	0.535%
79	14.420	2184	2187	2193	rVB2	45761	59990	3.13%	0.352%
80	14.487	2193	2198	2202	rBV2	91947	154105	8.05%	0.904%
81	14.578	2210	2213	2216	rBV3	23941	29395	1.54%	0.173%
82	14.615	2216	2219	2225	rVV5	70112	106126	5.54%	0.623%
83	14.676	2225	2229	2233	rVB	77071	97194	5.08%	0.570%
84	14.725	2233	2237	2241	rBV2	29836	42841	2.24%	0.251%
85	14.792	2243	2248	2251	rBV2	50891	80992	4.23%	0.475%
86	14.847	2254	2257	2261	rVB3	51944	65284	3.41%	0.383%
87	15.048	2283	2290	2293	rBV7	19290	41461	2.17%	0.243%
88	15.182	2307	2312	2320	rVV3	66255	132478	6.92%	0.778%
89	15.255	2320	2324	2330	rVB5	27017	46660	2.44%	0.274%
90	15.377	2340	2344	2346	rBV4	18894	27383	1.43%	0.161%
91	15.407	2346	2349	2353	rVV2	32192	46725	2.44%	0.274%
92	15.481	2353	2361	2364	rVV3	33358	57029	2.98%	0.335%
93	15.523	2365	2368	2374	rVV4	19120	37838	1.98%	0.222%
94	15.615	2378	2383	2386	rVV2	52023	85529	4.47%	0.502%
95	15.651	2386	2389	2397	rVB6	31496	55561	2.90%	0.326%
96	15.731	2397	2402	2409	rBV3	19696	43971	2.30%	0.258%
97	15.810	2410	2415	2417	rVV4	17988	30278	1.58%	0.178%
98	15.846	2417	2421	2429	rVB4	25034	50432	2.63%	0.296%

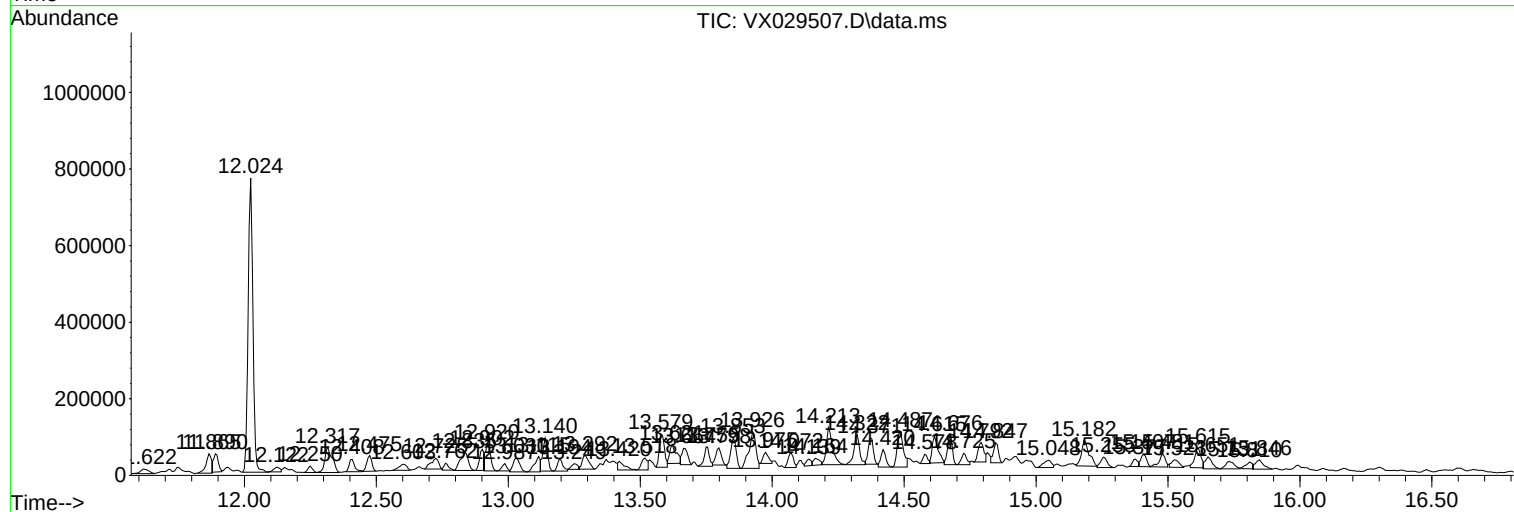
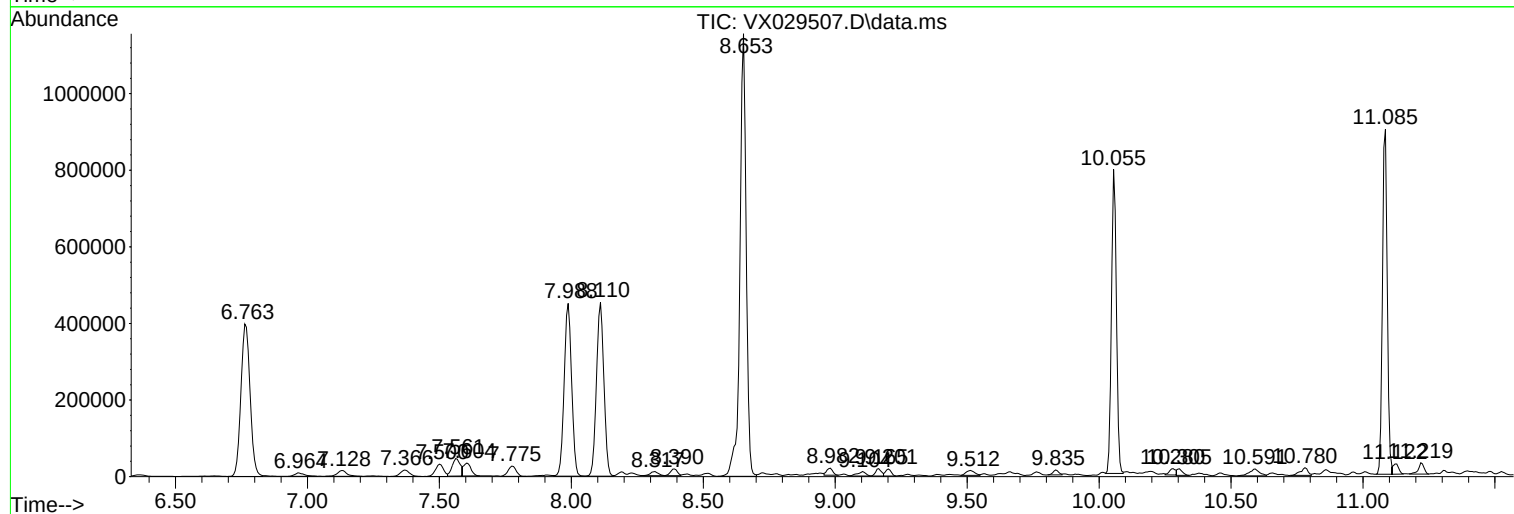
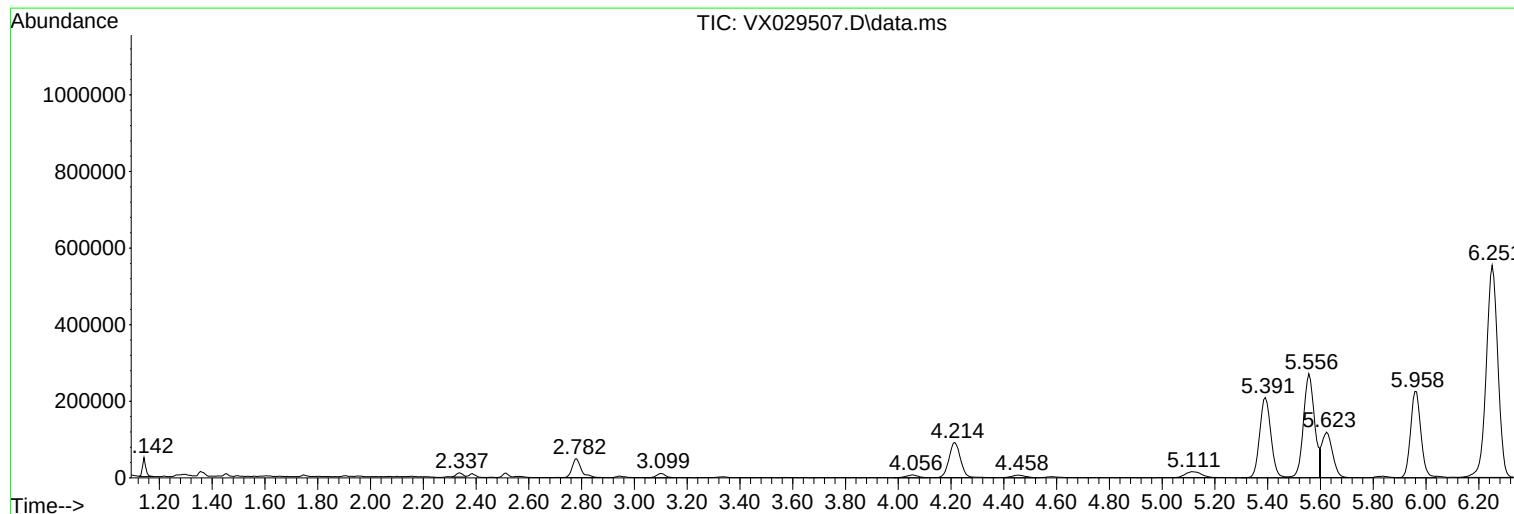
Sum of corrected areas: 17037955

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Instrument :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060722W.M
Quant Title : SW846 8260

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



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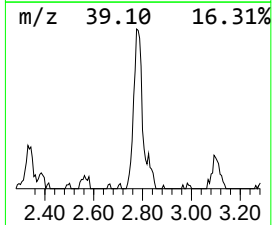
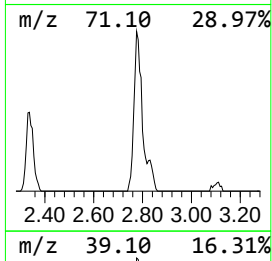
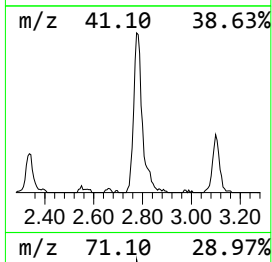
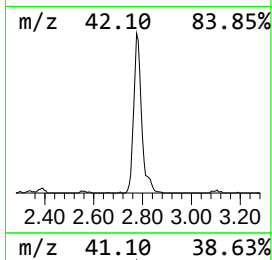
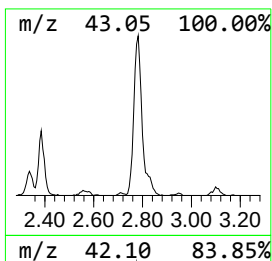
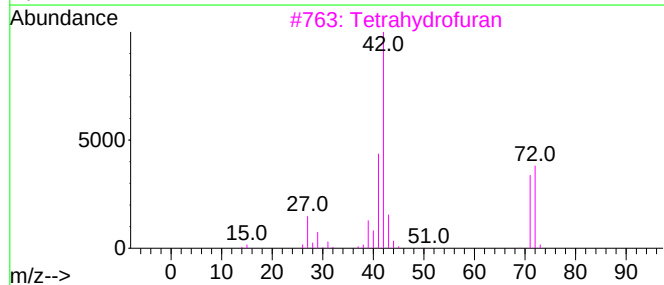
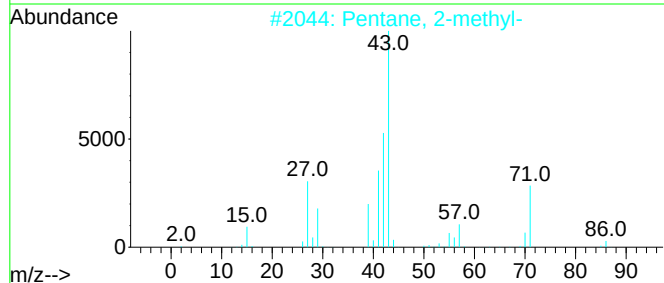
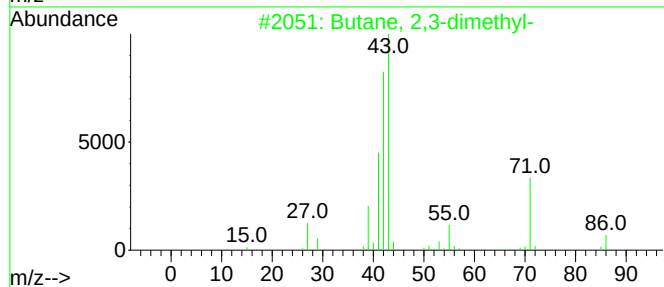
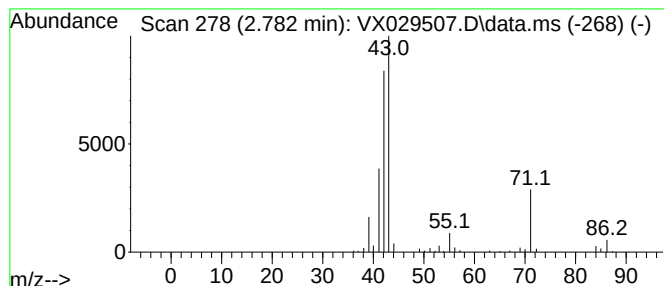
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060722W.M
Quant Title : SW846 8260

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

Peak Number 1 Butane, 2,3-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.782	7.39 ug/l	119392	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	91
2			Pentane, 2-methyl-	86	C6H14	000107-83-5	50
3			Tetrahydrofuran	72	C4H8O	000109-99-9	38
4			Pyrrolidine	71	C4H9N	000123-75-1	16
5			C2H5C(CH3)2COCH3	114	C7H14O	020669-04-9	12



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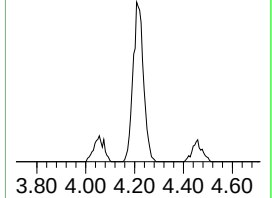
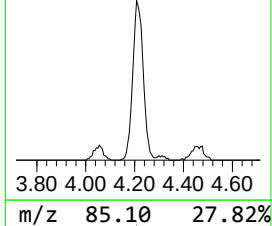
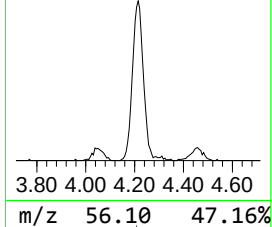
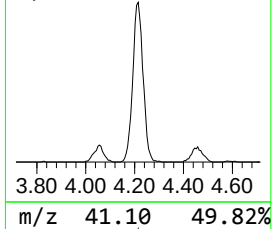
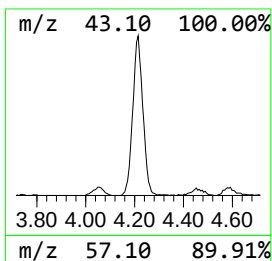
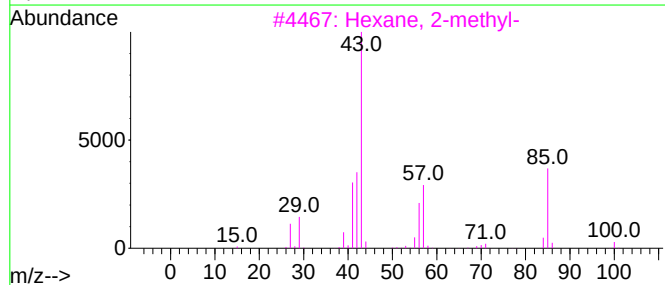
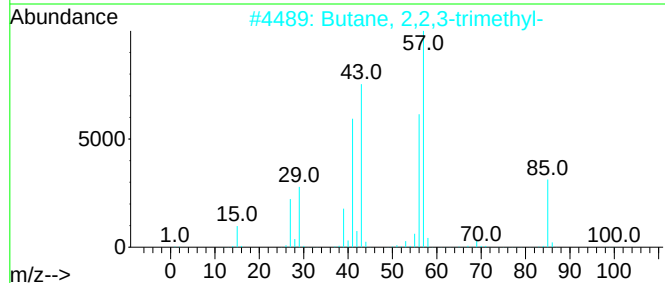
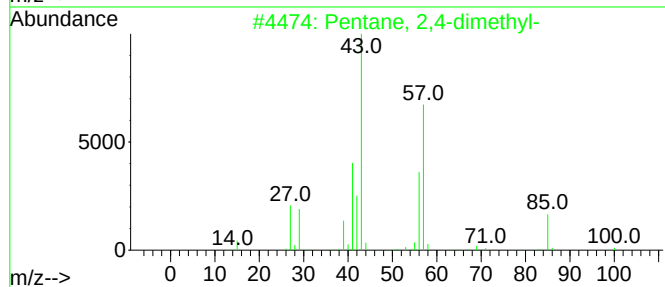
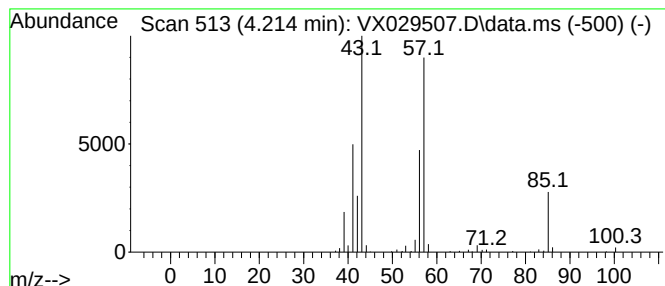
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060722W.M
Quant Title : SW846 8260

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

Peak Number 2 Pentane, 2,4-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.214	17.02 ug/l	275231	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	91
2			Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	64
3			Hexane, 2-methyl-	100	C7H16	000591-76-4	64
4			Oxalic acid, butyl isohexyl ester	230	C12H22O4	1000309-32-7	64
5			Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	59



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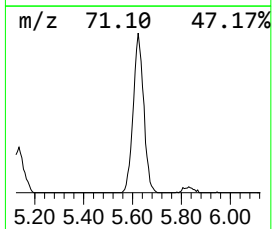
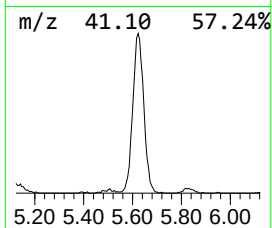
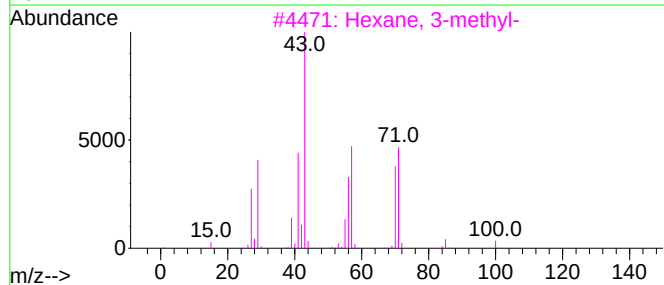
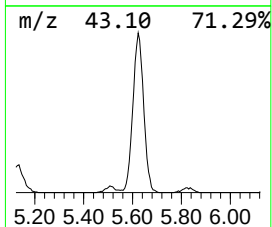
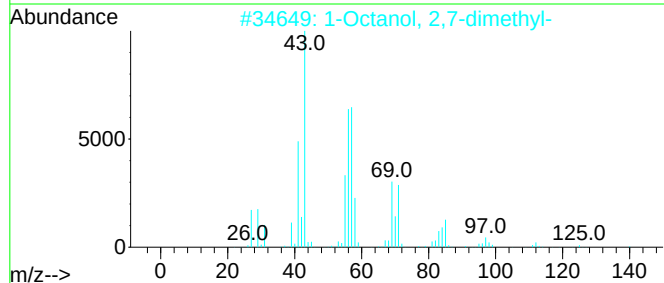
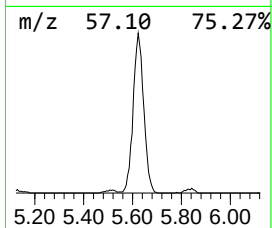
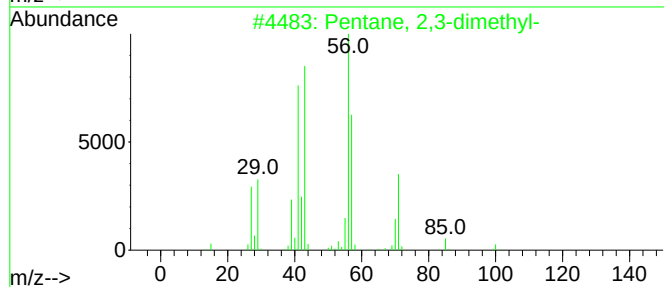
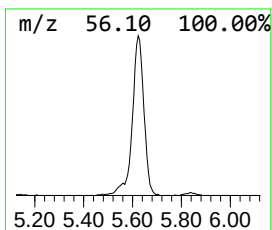
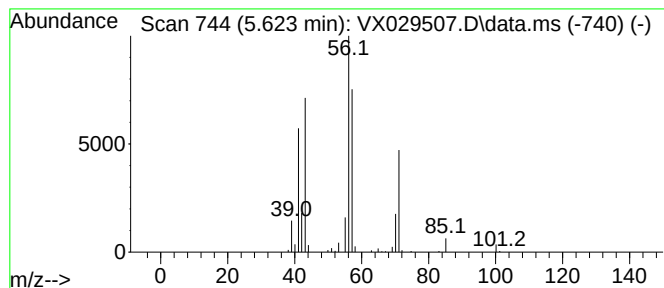
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060722W.M
Quant Title : SW846 8260

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

Peak Number 3 Pentane, 2,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.623	19.83 ug/l	320618	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
2			1-Octanol, 2,7-dimethyl-	158	C10H22O	015250-22-3	40
3			Hexane, 3-methyl-	100	C7H16	000589-34-4	38
4			Heptane, 2,2,4-trimethyl-	142	C10H22	014720-74-2	38
5			Octane	114	C8H18	000111-65-9	37



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061522\
Data File : VX029507.D
Acq On : 15 Jun 2022 14:38
Operator : JC/MD
Sample : N3160-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2

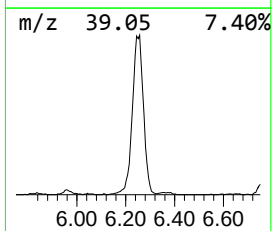
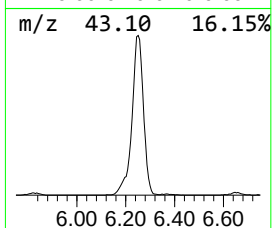
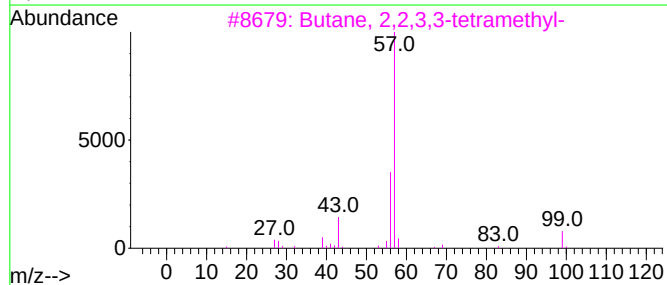
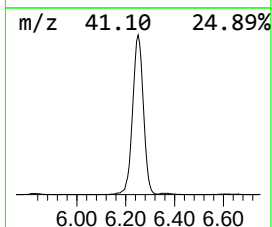
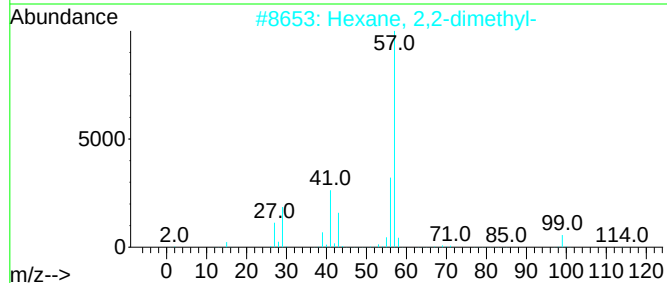
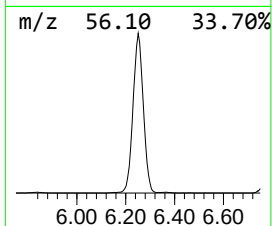
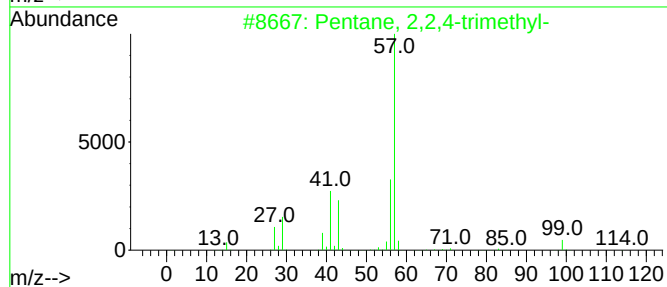
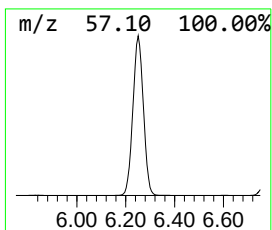
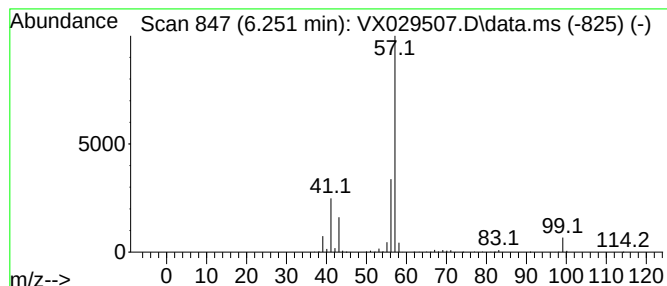
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Quant Title : SW846 8260

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

Peak Number 4 Pentane, 2,2,4-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.251	88.63 ug/l	1689350	1,4-Difluorobenzene	6.763

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	83
2			Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	83
3			Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	78
4			Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	72
5			Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061522\
Data File : VX029507.D
Acq On : 15 Jun 2022 14:38
Operator : JC/MD
Sample : N3160-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2

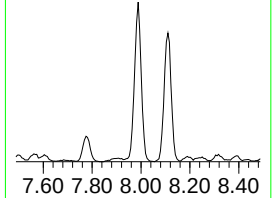
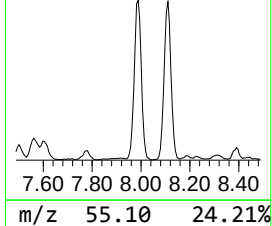
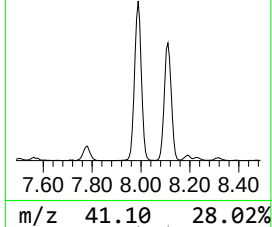
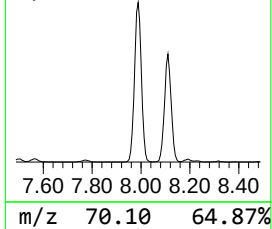
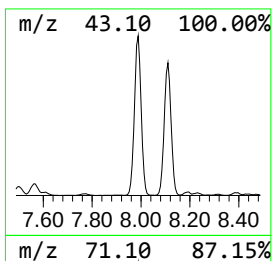
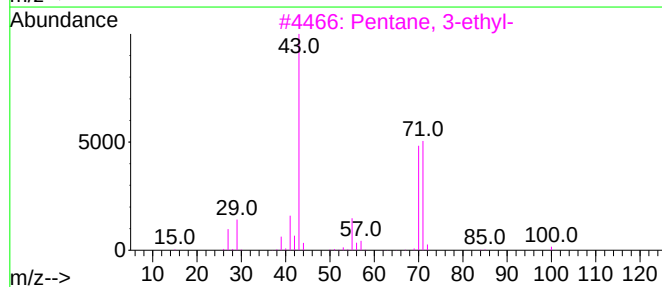
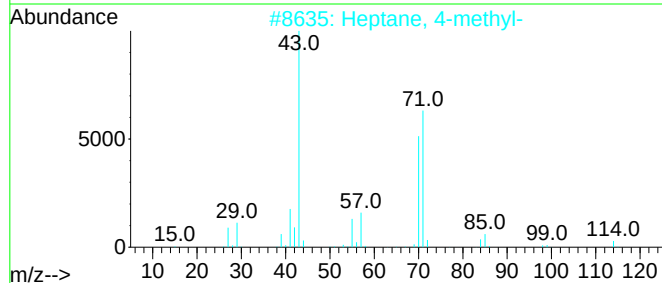
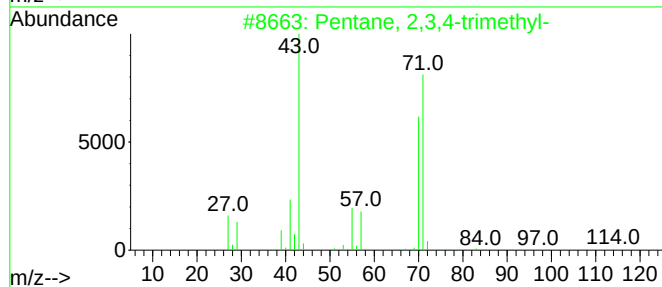
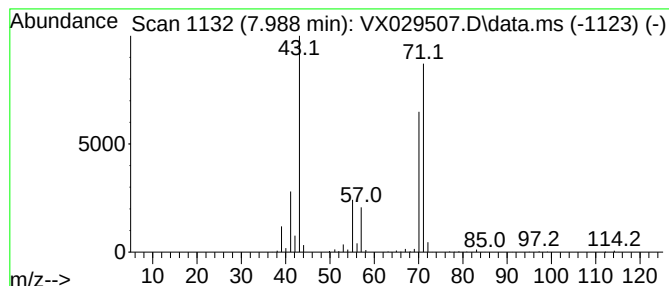
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Quant Title : SW846 8260

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

Peak Number 5 Pentane, 2,3,4-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.988	46.31 ug/l	882625	1,4-Difluorobenzene	6.763

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	91
2			Heptane, 4-methyl-	114	C8H18	000589-53-7	83
3			Pentane, 3-ethyl-	100	C7H16	000617-78-7	83
4			Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	78
5			Diisooamyl ether	158	C10H22O	000544-01-4	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061522\
Data File : VX029507.D
Acq On : 15 Jun 2022 14:38
Operator : JC/MD
Sample : N3160-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2

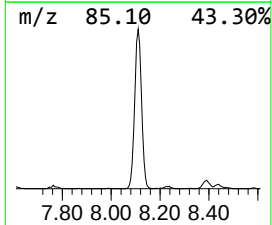
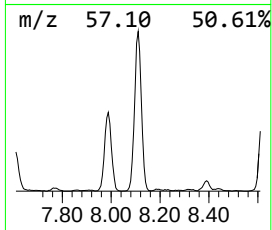
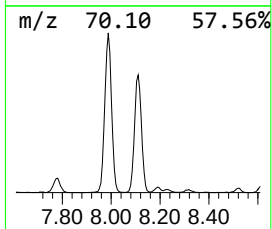
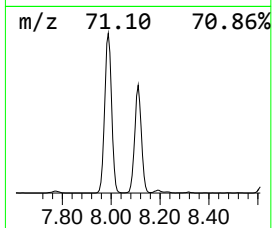
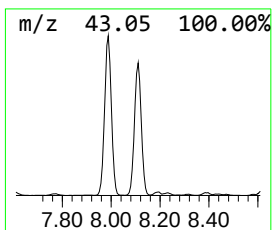
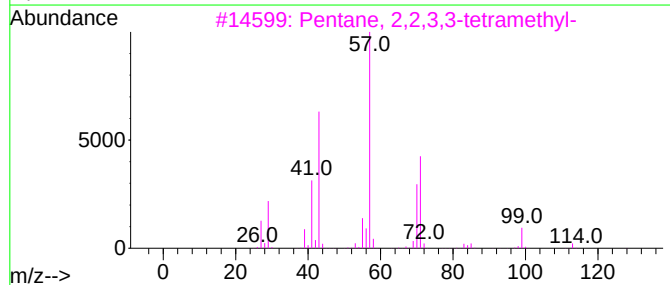
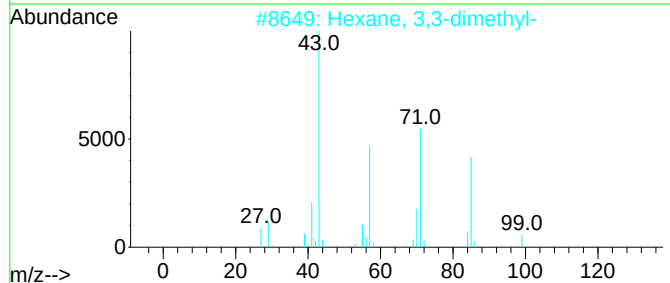
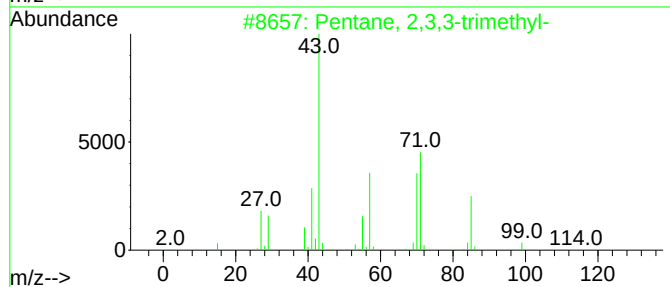
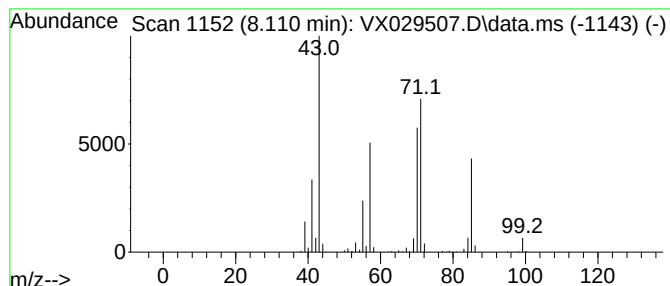
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Quant Title : SW846 8260

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

Peak Number 6 Pentane, 2,3,3-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.110	46.51 ug/l	886401	1,4-Difluorobenzene	6.763

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2		Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	72
3		Pentane, 2,2,3,3-tetramethyl-	128	C9H20	007154-79-2	53
4		Hexane, 3-methyl-	100	C7H16	000589-34-4	53
5		Heptane, 4-methyl-	114	C8H18	000589-53-7	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061522\
Data File : VX029507.D
Acq On : 15 Jun 2022 14:38
Operator : JC/MD
Sample : N3160-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

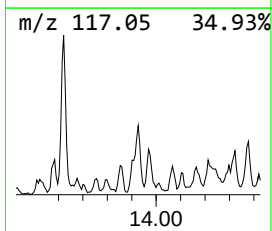
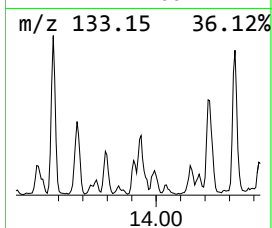
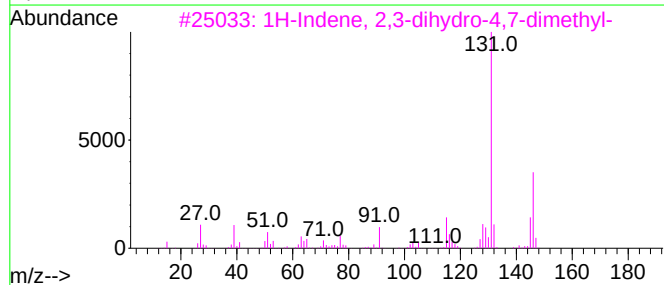
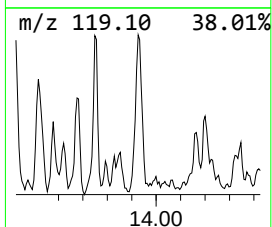
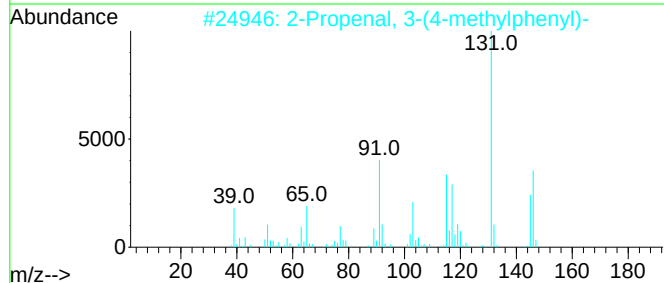
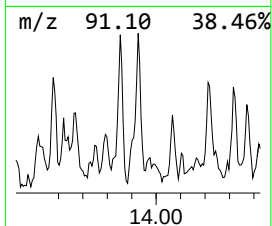
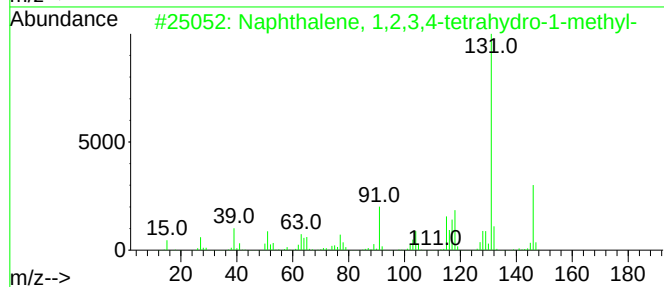
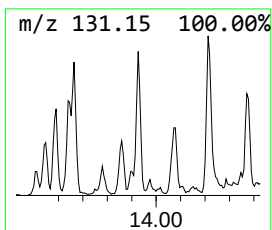
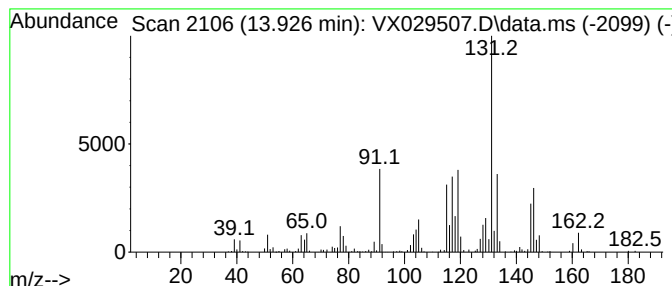
Instrument :
MSVOA_X
ClientSampleId :
MW-2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060722W.M
Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD			R.T.
13.926	9.18 ug/l	178323	1,4-Dichlorobenzene-d4			12.024
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...		146	C11H14	001559-81-5	92
2	2-Propenal, 3-(4-methylphenyl)-		146	C10H10O	001504-75-2	70
3	1H-Indene, 2,3-dihydro-4,7-dimet...		146	C11H14	006682-71-9	64
4	Benzene, (3-methyl-2-butenyl)-		146	C11H14	004489-84-3	64
5	Benzene, (2-methyl-1-butenyl)-		146	C11H14	056253-64-6	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061522\
Data File : VX029507.D
Acq On : 15 Jun 2022 14:38
Operator : JC/MD
Sample : N3160-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2

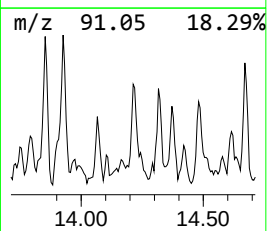
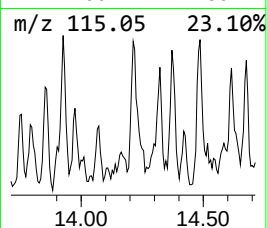
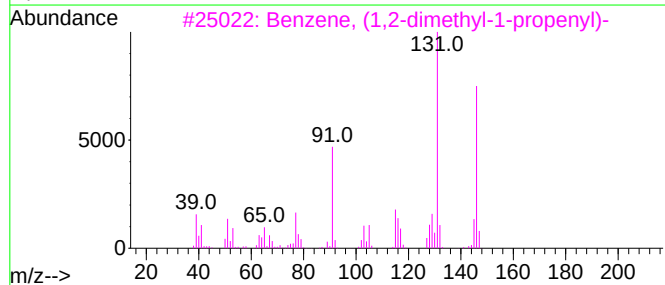
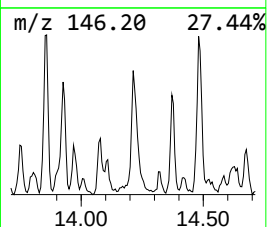
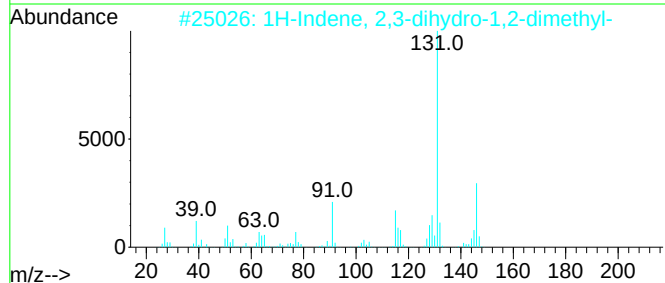
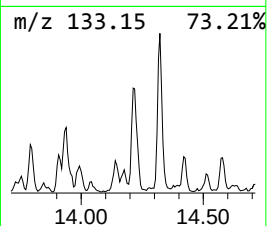
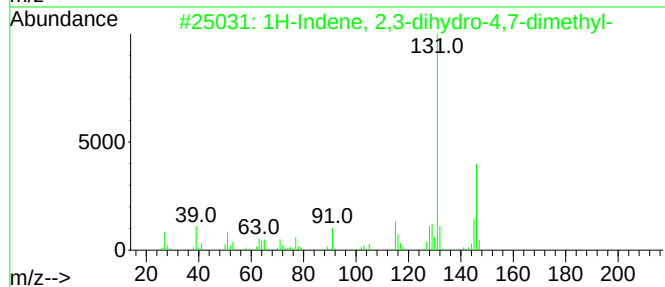
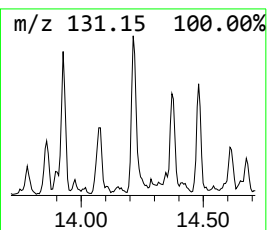
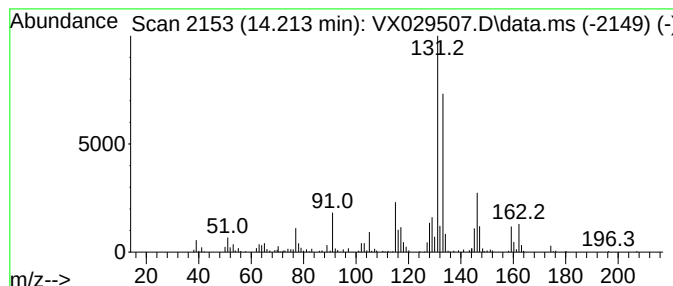
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Quant Title : SW846 8260

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

Peak Number 8 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.213	8.22 ug/l	159619	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	87
2			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	70
3			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	70
4			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	55
5			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061522\
Data File : VX029507.D
Acq On : 15 Jun 2022 14:38
Operator : JC/MD
Sample : N3160-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2

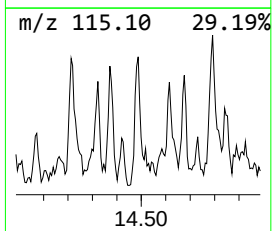
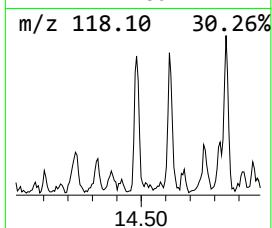
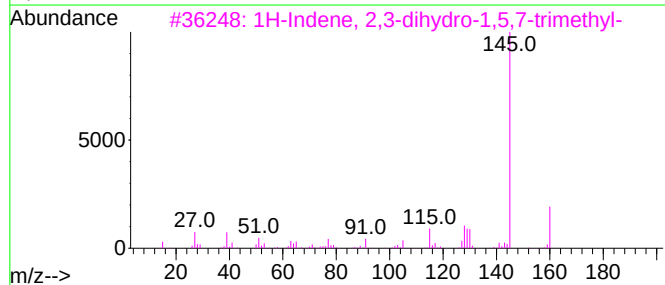
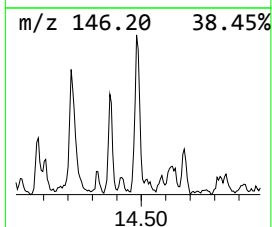
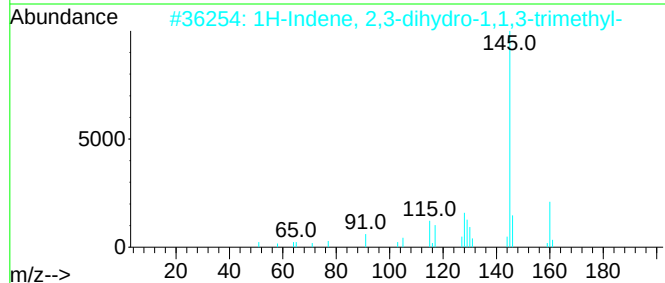
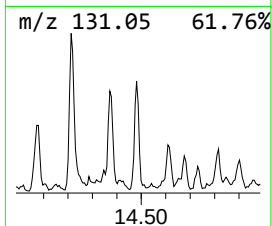
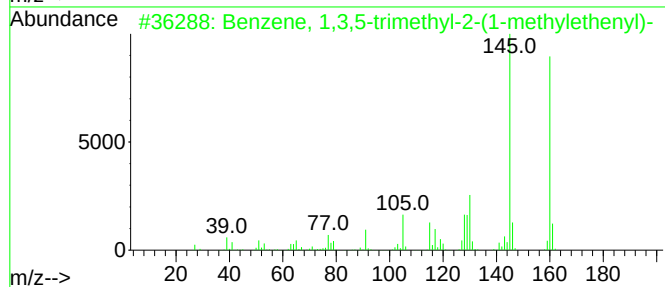
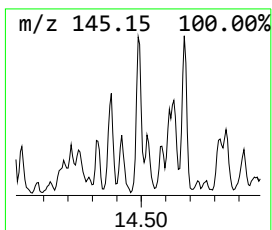
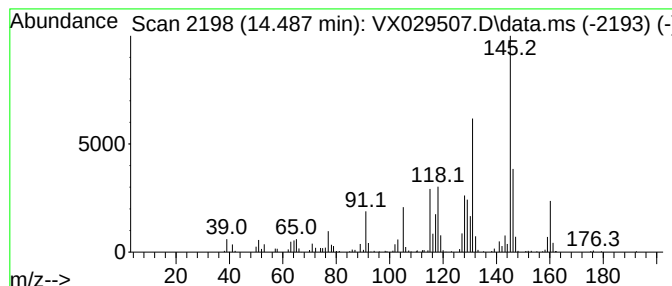
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,3,5-trimethyl-2-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.487	7.93 ug/l	154105	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	70
2			1H-Indene, 2,3-dihydro-1,1,3-tri...	160	C12H16	002613-76-5	64
3			1H-Indene, 2,3-dihydro-1,5,7-tri...	160	C12H16	054340-88-4	60
4			Benzene, 1-(1-methylethenyl)-2-(...	160	C12H16	005557-93-7	58
5			Benzene, 1-(1-methylethenyl)-3-(...	160	C12H16	001129-29-9	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061522\
Data File : VX029507.D
Acq On : 15 Jun 2022 14:38
Operator : JC/MD
Sample : N3160-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

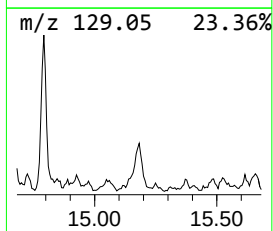
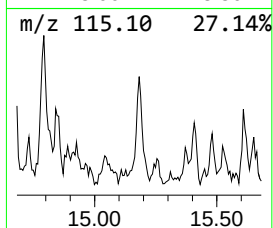
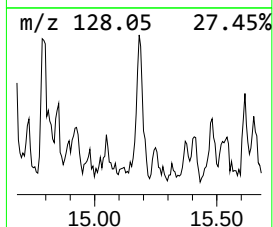
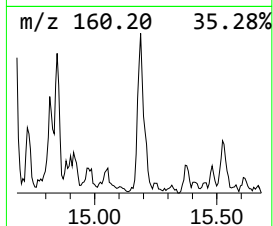
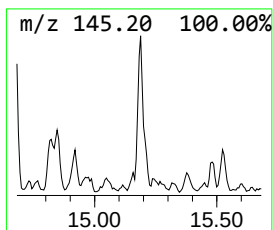
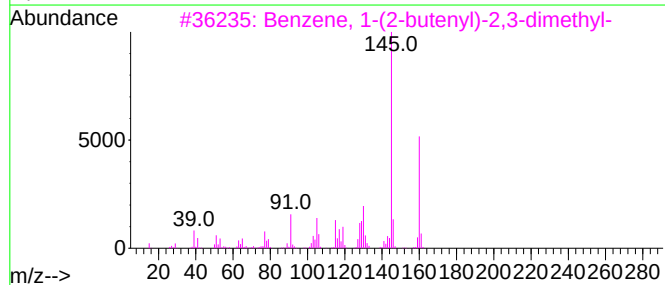
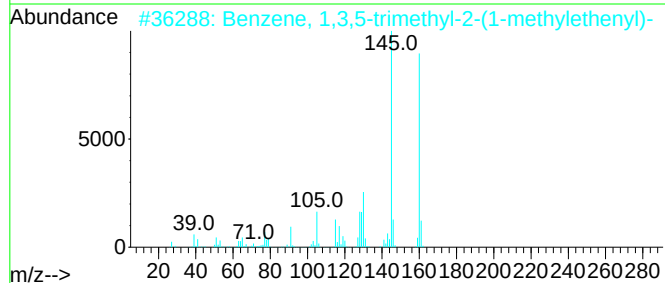
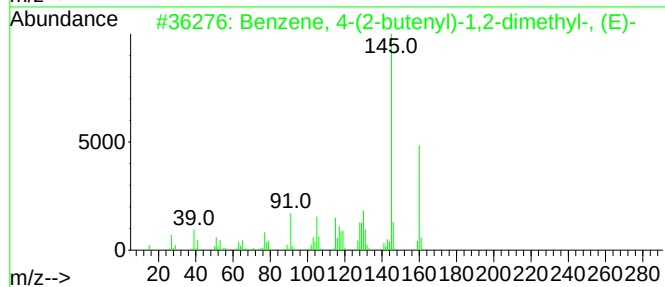
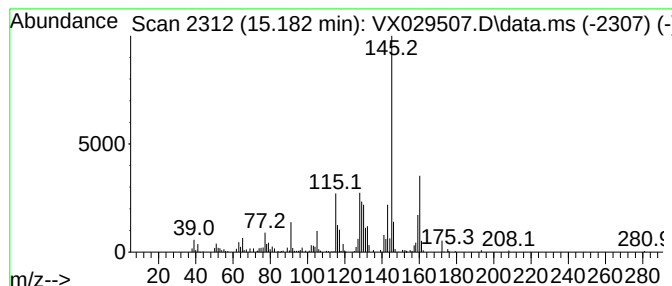
Instrument :
MSVOA_X
ClientSampleId :
MW-2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060722W.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, 4-(2-butenyl)-1,2-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD			R.T.
15.182	6.82 ug/l	132478	1,4-Dichlorobenzene-d4			12.024
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, 4-(2-butenyl)-1,2-dimet...		160	C12H16	054340-86-2	70
2	Benzene, 1,3,5-trimethyl-2-(1-me...		160	C12H16	014679-13-1	70
3	Benzene, 1-(2-butenyl)-2,3-dimet...		160	C12H16	054340-85-1	62
4	Naphthalene, 1,2,3,4-tetrahydro-...		160	C12H16	001076-61-5	58
5	Naphthalene, 1,2,3,4-tetrahydro-...		160	C12H16	004175-54-6	52



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061522\
Data File : VX029507.D
Acq On : 15 Jun 2022 14:38
Operator : JC/MD
Sample : N3160-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060722W.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, 2,3-dim...	2.782	7.4	ug/l	119392	1	5.556	808312	50.0
Pentane, 2,4-di...	4.214	17.0	ug/l	275231	1	5.556	808312	50.0
Pentane, 2,3-di...	5.623	19.8	ug/l	320618	1	5.556	808312	50.0
Pentane, 2,2,4-...	6.251	88.6	ug/l	1689350	2	6.763	953007	50.0
Pentane, 2,3,4-...	7.988	46.3	ug/l	882625	2	6.763	953007	50.0
Pentane, 2,3,3-...	8.110	46.5	ug/l	886401	2	6.763	953007	50.0
Naphthalene, 1,...	13.926	9.2	ug/l	178323	4	12.024	971154	50.0
1H-Indene, 2,3-...	14.213	8.2	ug/l	159619	4	12.024	971154	50.0
Benzene, 1,3,5-...	14.487	7.9	ug/l	154105	4	12.024	971154	50.0
Benzene, 4-(2-b...	15.182	6.8	ug/l	132478	4	12.024	971154	50.0