

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051324\
 Data File : VX041482.D
 Acq On : 13 May 2024 11:13
 Operator : JC/MD
 Sample : P2413-01ME
 Misc : 4.99g/10mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 WC-CCME

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\82X043024W.M
 Title : SW846 8260

Signal : TIC: VX041481.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.794	274	280	293	rVB	37819	84825	10.13%	0.547%
2	3.074	317	326	340	rVB	35397	81109	9.68%	0.523%
3	4.184	499	508	516	rBV2	12066	35802	4.27%	0.231%
4	4.282	516	524	536	rVB3	17531	50115	5.98%	0.323%
5	5.385	691	705	713	rBV2	85436	274675	32.80%	1.772%
6	5.489	713	722	725	rVV4	76049	251318	30.01%	1.621%
7	5.544	725	731	762	rVB	169603	619609	73.98%	3.997%
8	5.806	762	774	789	rBV	102434	309786	36.99%	1.999%
9	5.952	789	798	816	rBV	92466	251897	30.08%	1.625%
10	6.147	819	830	838	rVV3	36748	128968	15.40%	0.832%
11	6.239	838	845	854	rVV4	50088	149520	17.85%	0.965%
12	6.342	854	862	874	rVB2	32403	93243	11.13%	0.602%
13	6.757	921	930	939	rBV	236701	602809	71.97%	3.889%
14	7.165	987	997	1002	rBV2	15811	37686	4.50%	0.243%
15	7.366	1018	1030	1040	rBV2	167445	432522	51.64%	2.790%
16	7.488	1042	1050	1055	rBV	36429	74398	8.88%	0.480%
17	7.549	1055	1060	1070	rVB	47235	96870	11.57%	0.625%
18	7.647	1070	1076	1086	rVB	32304	65433	7.81%	0.422%
19	7.763	1086	1095	1103	rVB2	39981	85605	10.22%	0.552%
20	7.976	1120	1130	1142	rVB4	29880	103223	12.32%	0.666%
21	8.098	1142	1150	1153	rBV3	38002	76198	9.10%	0.492%
22	8.135	1153	1156	1160	rVV	26504	40217	4.80%	0.259%
23	8.177	1160	1163	1168	rVB2	36454	57080	6.82%	0.368%
24	8.269	1173	1178	1181	rBV	81504	134394	16.05%	0.867%
25	8.305	1181	1184	1191	rVB2	73763	126256	15.07%	0.815%
26	8.378	1191	1196	1200	rBV2	17062	30288	3.62%	0.195%
27	8.433	1200	1205	1209	rBV	133425	222041	26.51%	1.432%
28	8.500	1213	1216	1225	rVB3	34407	70482	8.42%	0.455%
29	8.598	1225	1232	1235	rBV3	85305	169662	20.26%	1.095%
30	8.647	1235	1240	1250	rVB	503848	837528	100.00%	5.403%
31	8.769	1250	1260	1264	rBV2	44157	83802	10.01%	0.541%
32	8.811	1264	1267	1270	rBV	26776	38843	4.64%	0.251%
33	8.915	1277	1284	1290	rBV4	39720	84113	10.04%	0.543%
34	8.970	1290	1293	1299	rVV	34511	57801	6.90%	0.373%
35	9.098	1306	1314	1319	rBV	53531	101770	12.15%	0.657%
36	9.159	1319	1324	1328	rVV	45249	72552	8.66%	0.468%

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Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X043024W.M
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37	9.195	1328	1330	1340	rVB5	18986	31791	3.80%	0.205%
38	9.378	1355	1360	1365	rBV3	32632	58144	6.94%	0.375%
39	9.494	1375	1379	1386	rVV3	63352	132754	15.85%	0.856%
40	9.561	1386	1390	1394	rVB	40261	55165	6.59%	0.356%
41	9.604	1394	1397	1402	rBV5	17734	37030	4.42%	0.239%
42	9.762	1416	1423	1428	rBV3	24601	48342	5.77%	0.312%
43	9.817	1428	1432	1436	rVV3	25740	54101	6.46%	0.349%
44	9.866	1436	1440	1442	rVV2	26525	48535	5.80%	0.313%
45	9.902	1442	1446	1452	rVV	115858	173848	20.76%	1.122%
46	10.006	1455	1463	1466	rVV	81748	117592	14.04%	0.759%
47	10.055	1466	1471	1477	rVV	518586	737306	88.03%	4.757%
48	10.128	1480	1483	1485	rVV3	18879	29813	3.56%	0.192%
49	10.195	1489	1494	1499	rVV	47907	87449	10.44%	0.564%
50	10.274	1499	1507	1510	rVV4	28926	77895	9.30%	0.503%
51	10.311	1510	1513	1517	rVV2	23598	40311	4.81%	0.260%
52	10.360	1517	1521	1533	rVB7	30914	73654	8.79%	0.475%
53	10.585	1552	1558	1561	rVV	39442	58260	6.96%	0.376%
54	10.671	1566	1572	1577	rBV3	22123	40891	4.88%	0.264%
55	10.774	1582	1589	1593	rBV2	30827	60862	7.27%	0.393%
56	10.854	1598	1602	1607	rVB3	23498	37816	4.52%	0.244%
57	10.963	1614	1620	1627	rBV	93963	131562	15.71%	0.849%
58	11.030	1627	1631	1634	rVV	19231	28640	3.42%	0.185%
59	11.079	1634	1639	1651	rVV	469499	716218	85.52%	4.621%
60	11.189	1651	1657	1665	rVV2	40144	64556	7.71%	0.416%
61	11.305	1671	1676	1684	rVV	355881	469482	56.06%	3.029%
62	11.402	1685	1692	1695	rVV	23592	49429	5.90%	0.319%
63	11.451	1695	1700	1711	rVB	101109	166046	19.83%	1.071%
64	11.616	1722	1727	1734	rBV	55030	87216	10.41%	0.563%
65	11.750	1746	1749	1760	rVB	80027	117260	14.00%	0.756%
66	11.866	1765	1768	1770	rVV	39565	49508	5.91%	0.319%
67	11.890	1770	1772	1776	rVV	44255	47115	5.63%	0.304%
68	12.024	1788	1794	1800	rVV	588489	782473	93.43%	5.048%
69	12.085	1800	1804	1811	rVV	60631	101134	12.08%	0.652%
70	12.244	1825	1830	1837	rBV2	385461	531198	63.42%	3.427%
71	12.317	1837	1842	1849	rVB3	240489	431081	51.47%	2.781%
72	12.402	1851	1856	1862	rBV2	27332	50324	6.01%	0.325%
73	12.463	1862	1866	1873	rVB	78433	93462	11.16%	0.603%
74	12.530	1873	1877	1879	rBV	64594	78433	9.36%	0.506%
75	12.554	1879	1881	1886	rVV	50286	59291	7.08%	0.383%
76	12.609	1886	1890	1894	rVV	389573	502636	60.01%	3.243%

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 Title : SW846 8260

77	12.646	1894	1896	1900	rVV	81030	85879	10.25%	0.554%
78	12.701	1900	1905	1913	rVV2	207627	339112	40.49%	2.188%
79	12.847	1925	1929	1933	rVB2	28292	37138	4.43%	0.240%
80	12.920	1933	1941	1945	rBV	207942	277728	33.16%	1.792%
81	12.963	1945	1948	1953	rVB	119583	146992	17.55%	0.948%
82	13.024	1953	1958	1960	rBV2	47108	66462	7.94%	0.429%
83	13.140	1974	1977	1978	rVV2	25631	32952	3.93%	0.213%
84	13.170	1978	1982	1986	rVV	191600	259107	30.94%	1.672%
85	13.213	1986	1989	1992	rVV	27114	33092	3.95%	0.213%
86	13.256	1992	1996	1998	rVV2	39476	58887	7.03%	0.380%
87	13.292	1998	2002	2011	rVV2	340405	498381	59.51%	3.215%
88	13.371	2011	2015	2020	rVV	43300	59673	7.12%	0.385%
89	13.420	2020	2023	2032	rVB2	34657	57161	6.82%	0.369%
90	13.506	2032	2037	2040	rBV2	34656	51291	6.12%	0.331%
91	13.542	2040	2043	2046	rVV	68837	87938	10.50%	0.567%
92	13.579	2046	2049	2055	rVV3	76948	139070	16.60%	0.897%
93	13.664	2056	2063	2069	rVV2	61158	125862	15.03%	0.812%
94	13.774	2076	2081	2088	rVV	107059	149688	17.87%	0.966%
95	13.853	2088	2094	2099	rVB5	13650	30697	3.67%	0.198%
96	13.926	2099	2106	2110	rBV3	31190	53867	6.43%	0.348%
97	14.072	2126	2130	2136	rBV	39391	49620	5.92%	0.320%
98	14.213	2148	2153	2163	rBV2	36030	67154	8.02%	0.433%
99	14.633	2208	2222	2232	rBV	90910	139277	16.63%	0.899%
100	14.780	2241	2246	2256	rVB	47896	62282	7.44%	0.402%

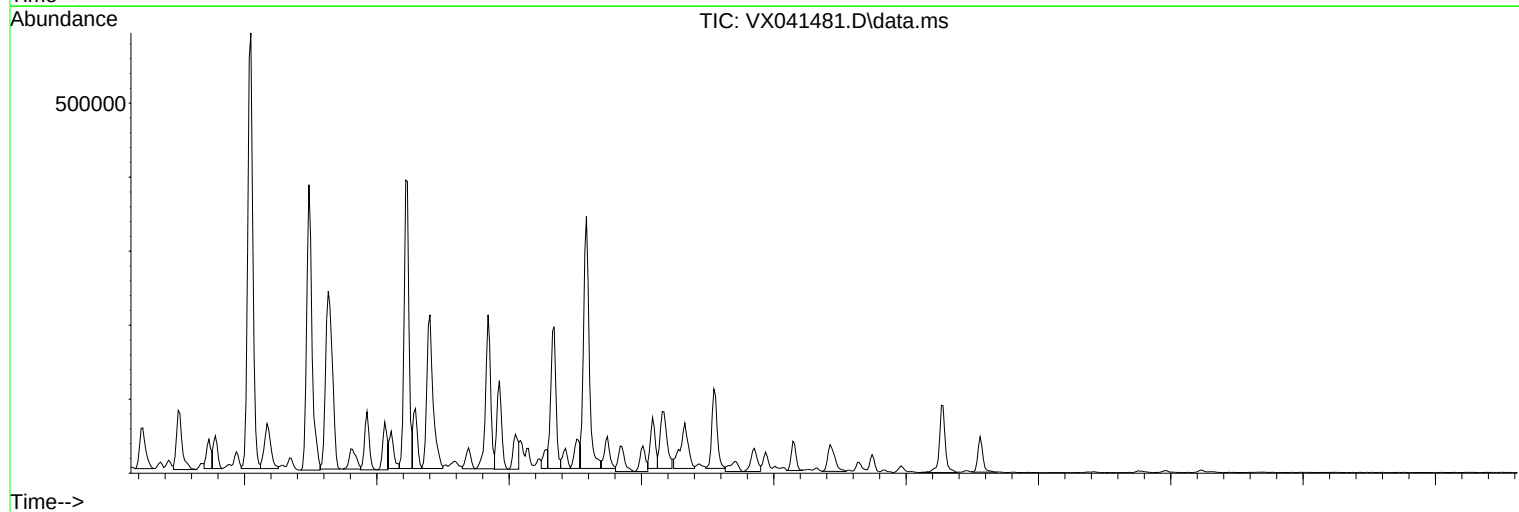
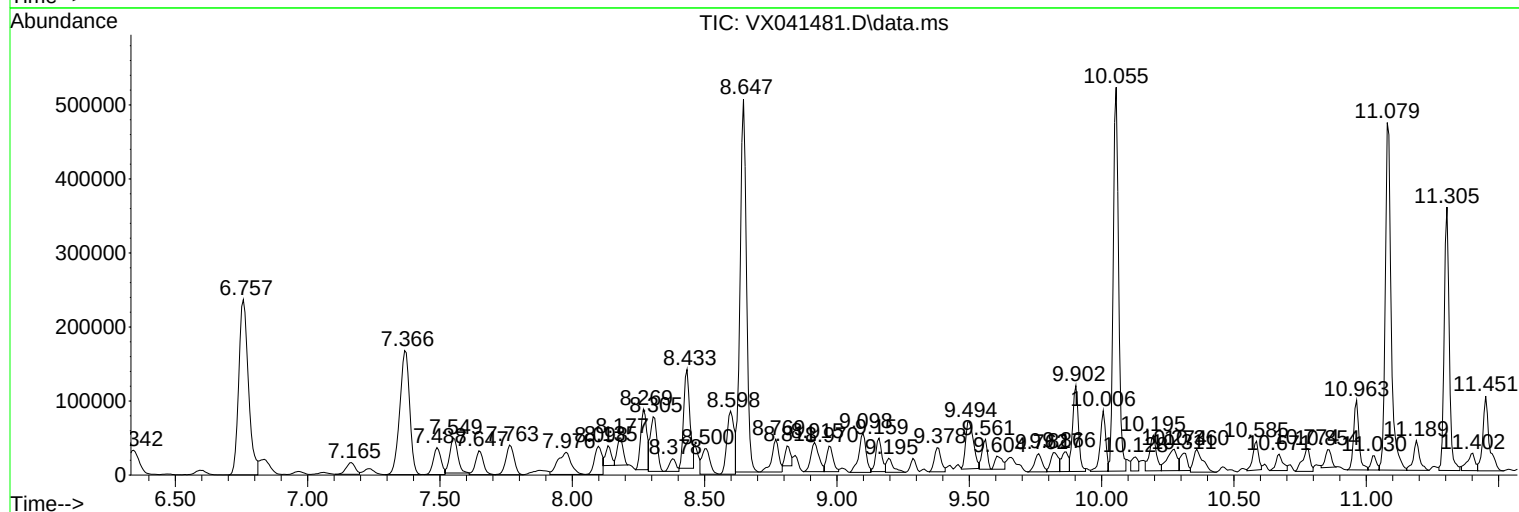
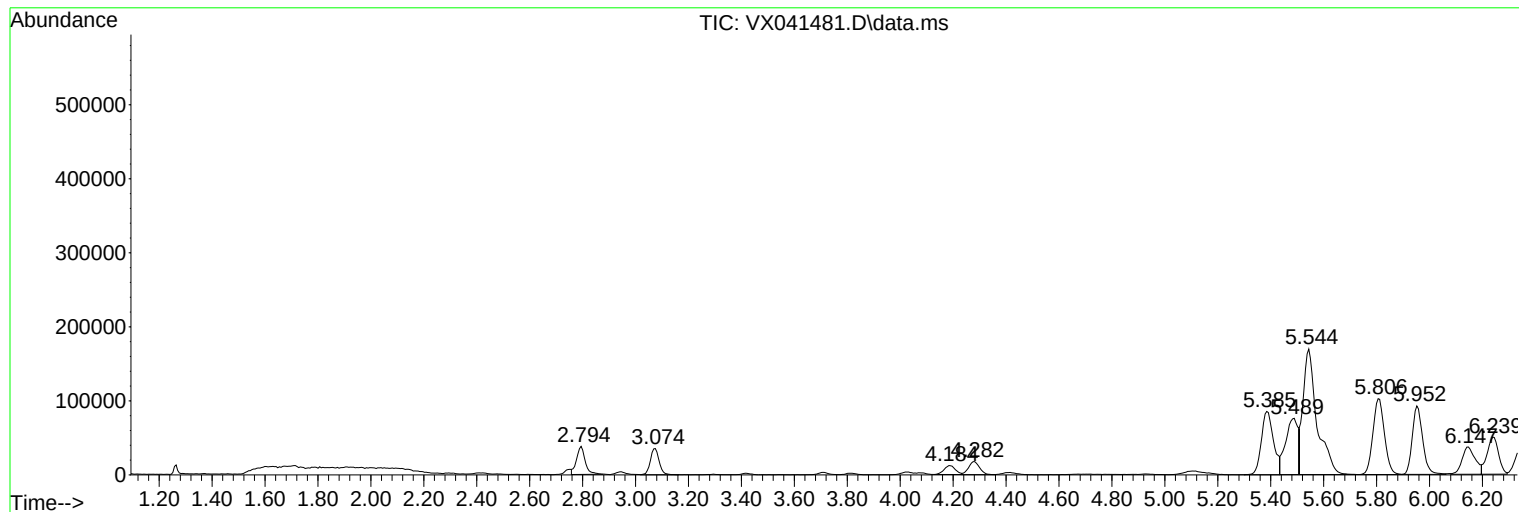
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Instrument :
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ClientSampleId :
WC-CCME

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X043024W.M
Quant Title : SW846 8260

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



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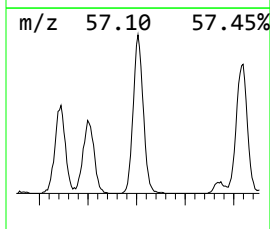
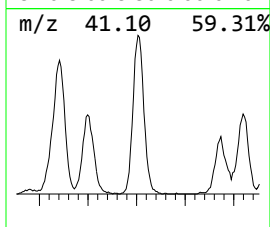
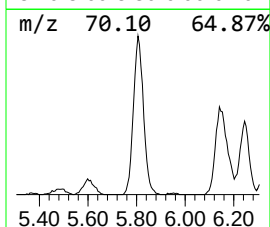
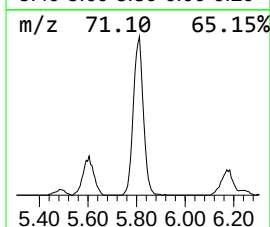
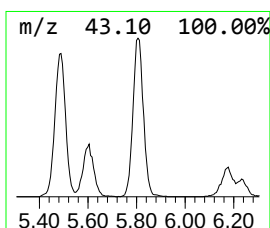
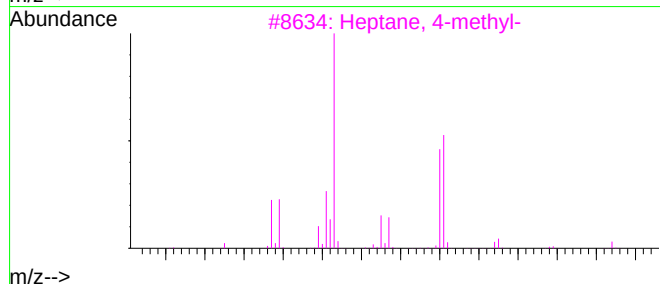
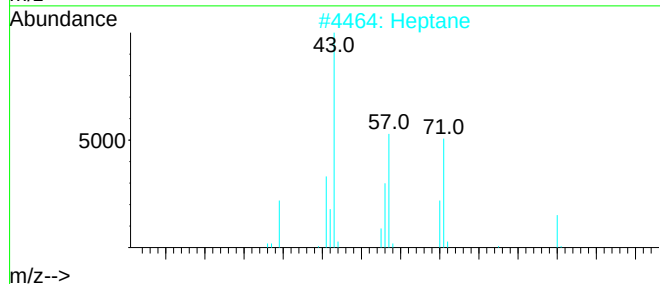
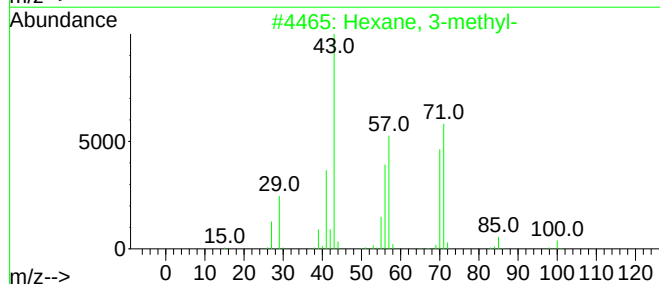
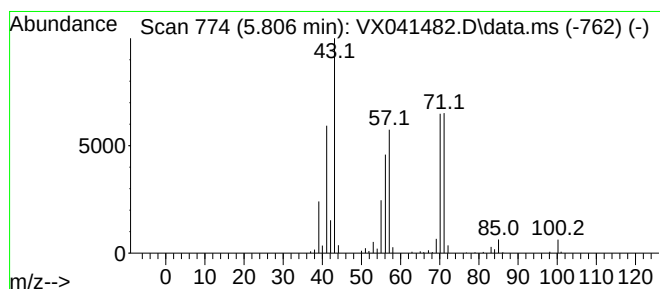
Instrument :
MSVOA_X
ClientSampleId :
WC-CCME

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X043024W.M
Quant Title : SW846 8260

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

Peak Number 1 Hexane, 3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.806	25.00 ug/l	309786	Pentafluorobenzene	5.544	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 3-methyl-	100	C7H16	000589-34-4	91
2	Heptane	100	C7H16	000142-82-5	80
3	Heptane, 4-methyl-	114	C8H18	000589-53-7	59
4	Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	53
5	1-Butanol, 2-ethyl-	102	C6H14O	000097-95-0	53



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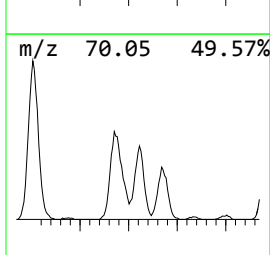
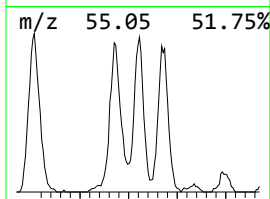
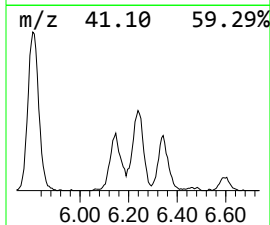
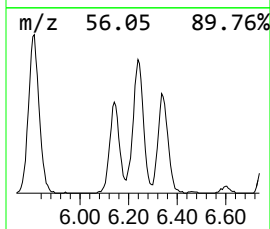
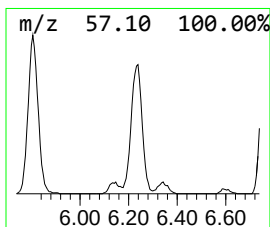
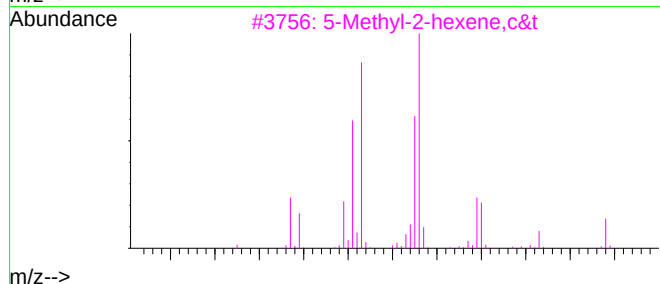
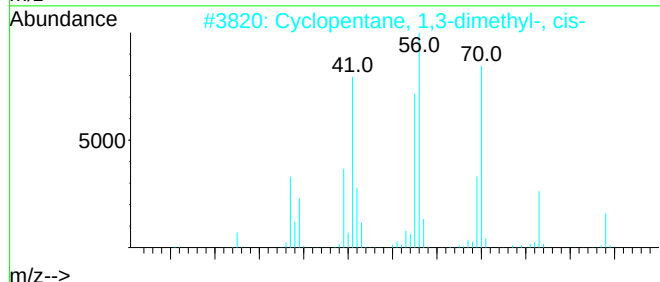
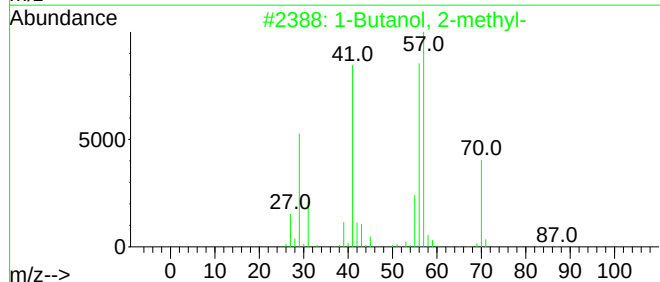
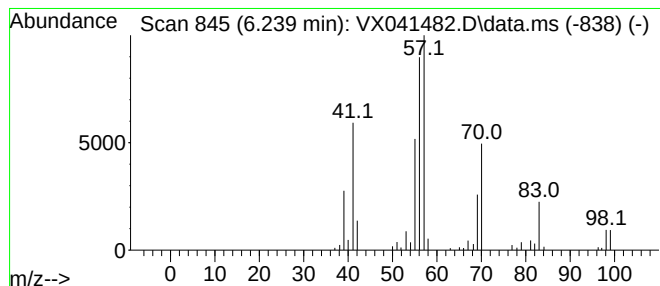
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X043024W.M
Quant Title : SW846 8260

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

Peak Number 2 1-Butanol, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.239	12.40 ug/l	149520	1,4-Difluorobenzene	6.757

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Butanol, 2-methyl-	88	C5H12O	000137-32-6	59
2	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	47
3	5-Methyl-2-hexene,c&t	98	C7H14	003404-62-4	47
4	1-Pentene, 2,4-dimethyl-	98	C7H14	002213-32-3	43
5	1-Hexene, 2-methyl-	98	C7H14	006094-02-6	43



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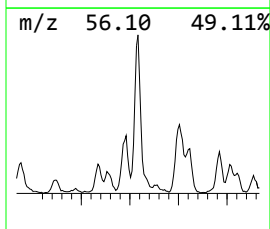
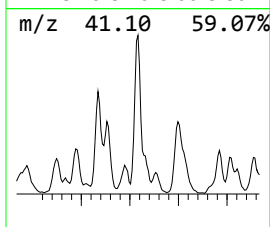
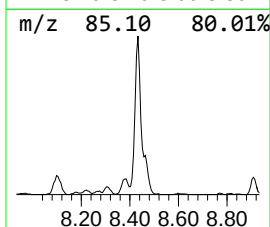
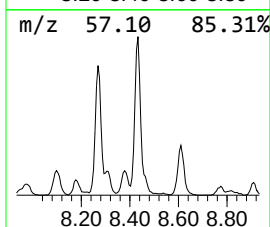
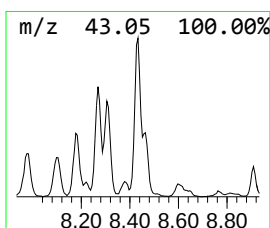
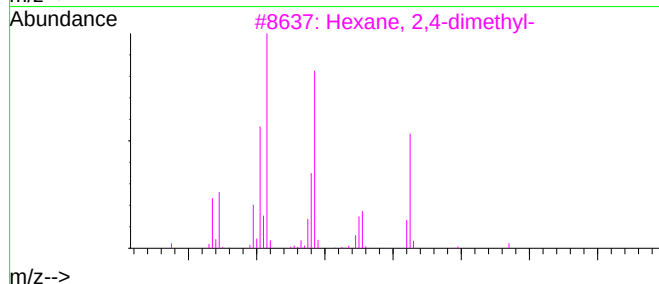
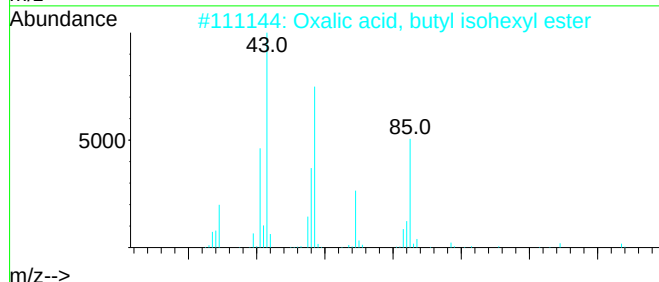
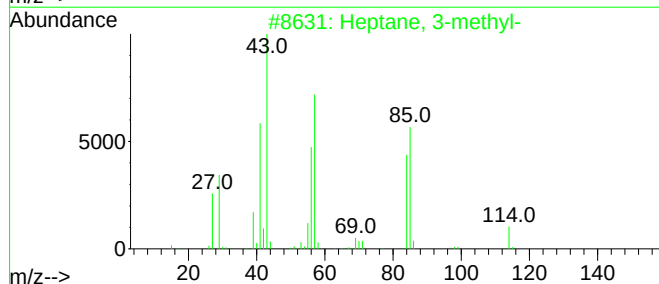
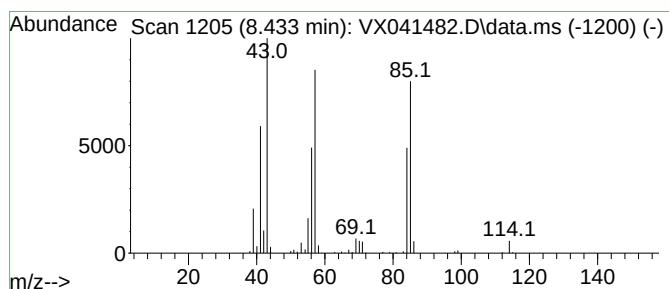
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Quant Title : SW846 8260

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

Peak Number 3 Heptane, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.433	15.06 ug/l	222041	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 3-methyl-	114	C8H18	000589-81-1	94
2	Oxalic acid, butyl isohexyl ester	230	C12H22O4	1000309-32-7	64
3	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	64
4	Hexane, 1-(hexyloxy)-2-methyl-	200	C13H28O	074421-17-3	64
5	Pentane, 3-ethyl-2,4-dimethyl-	128	C9H20	001068-87-7	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051324\
Data File : VX041482.D
Acq On : 13 May 2024 11:13
Operator : JC/MD
Sample : P2413-01ME
Misc : 4.99g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WC-CCME

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X043024W.M
Quant Title : SW846 8260

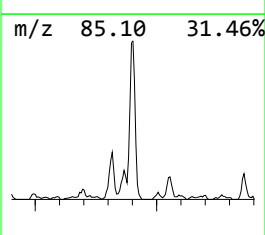
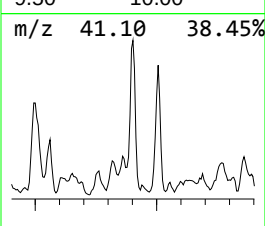
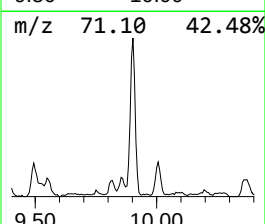
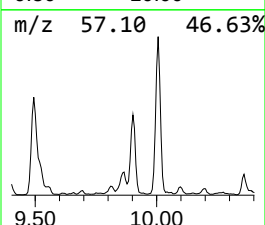
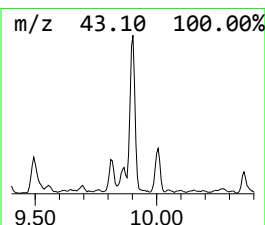
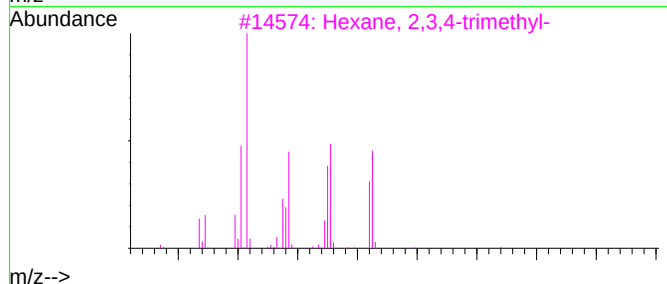
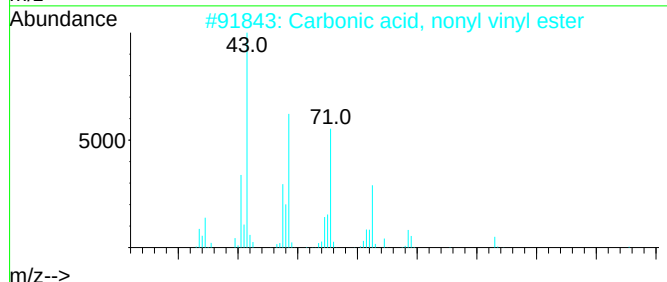
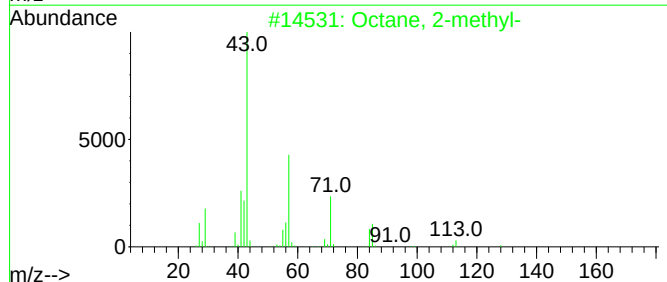
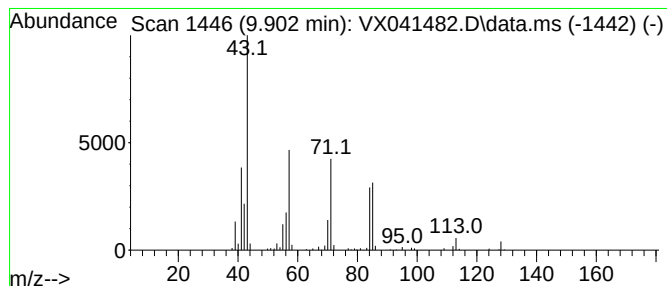
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 4 Octane, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.902	11.79 ug/l	173848	Chlorobenzene-d5	10.055

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2-methyl-	128	C9H20	003221-61-2	80
2			Carbonic acid, nonyl vinyl ester	214	C12H22O3	1000383-25-6	59
3			Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	59
4			Octane	114	C8H18	000111-65-9	58
5			Undecane, 4,7-dimethyl-	184	C13H28	017301-32-5	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051324\
Data File : VX041482.D
Acq On : 13 May 2024 11:13
Operator : JC/MD
Sample : P2413-01ME
Misc : 4.99g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WC-CCME

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X043024W.M
Quant Title : SW846 8260

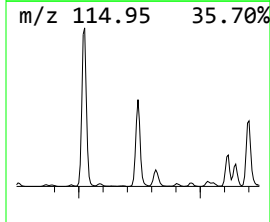
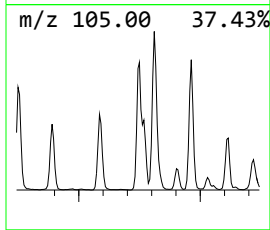
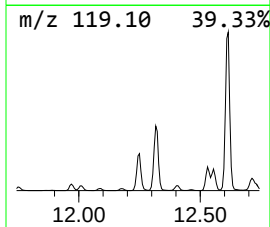
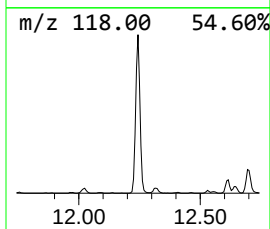
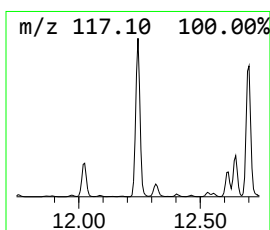
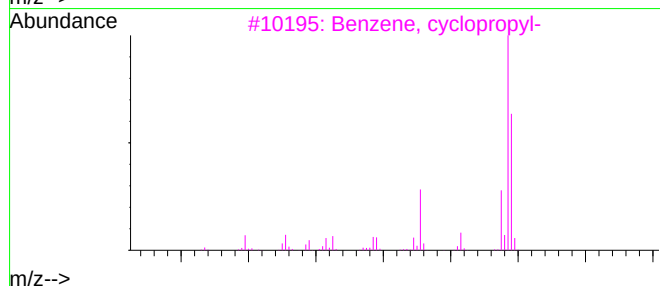
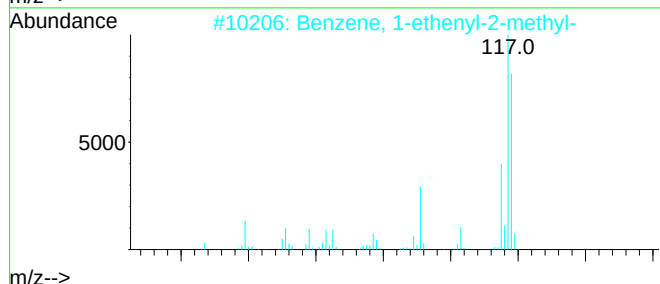
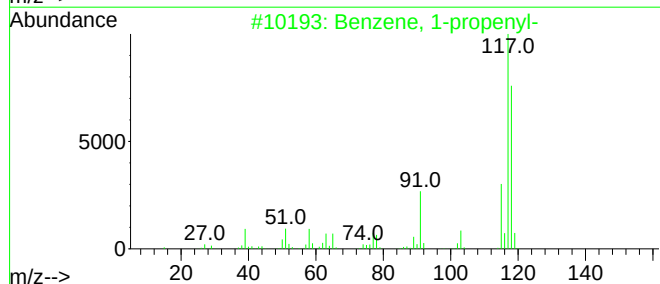
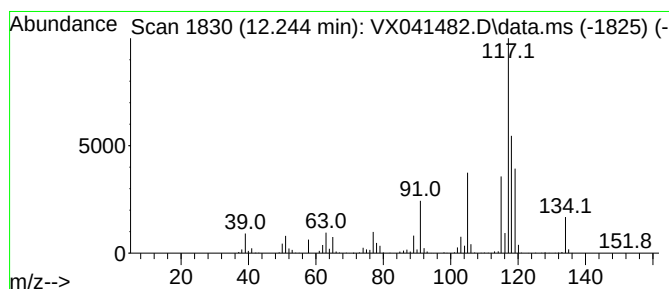
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TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1-propenyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	33.94 ug/l	531198	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-propenyl-	118	C9H10	000637-50-3	87
2	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	80
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	55
4	Indane	118	C9H10	000496-11-7	55
5	Benzene, 2-propenyl-	118	C9H10	000300-57-2	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051324\
Data File : VX041482.D
Acq On : 13 May 2024 11:13
Operator : JC/MD
Sample : P2413-01ME
Misc : 4.99g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WC-CCME

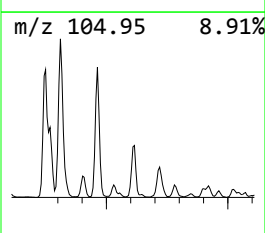
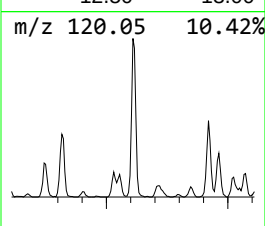
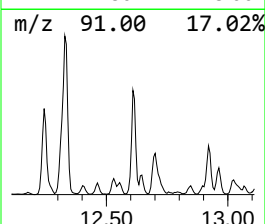
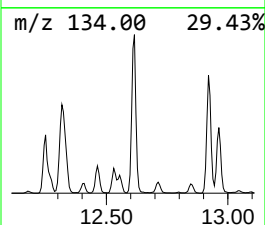
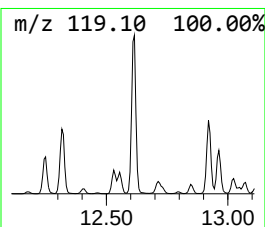
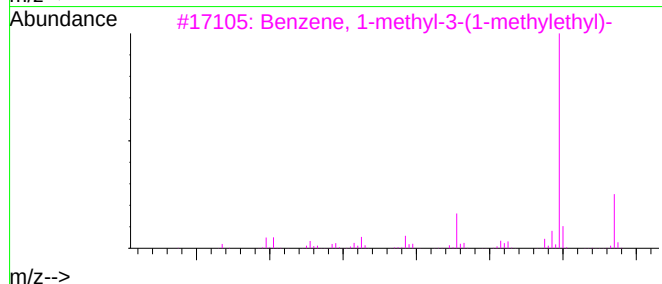
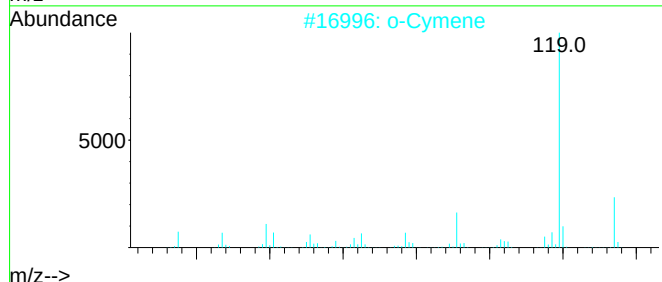
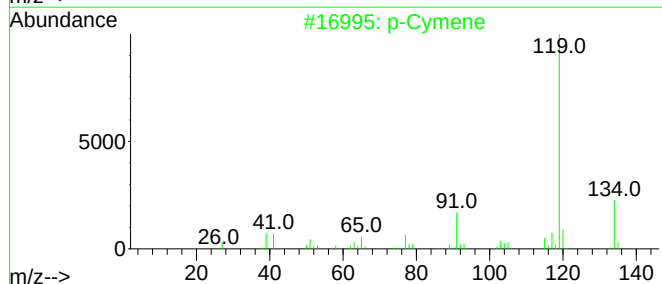
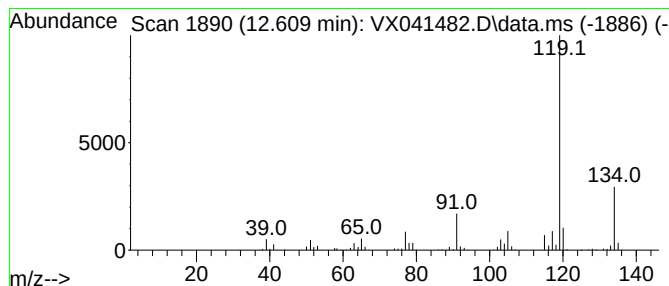
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 6 p-Cymene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.609	32.12 ug/l	502636	1,4-Dichlorobenzene-d4	12.024	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene	134	C10H14	000099-87-6	97
2	o-Cymene	134	C10H14	000527-84-4	97
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	97
4	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051324\
Data File : VX041482.D
Acq On : 13 May 2024 11:13
Operator : JC/MD
Sample : P2413-01ME
Misc : 4.99g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WC-CCME

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X043024W.M
Quant Title : SW846 8260

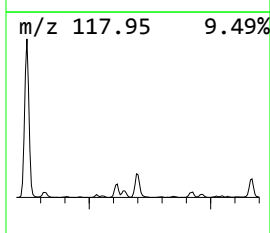
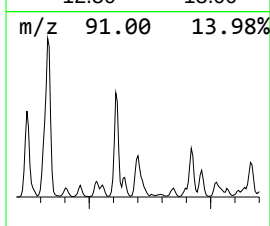
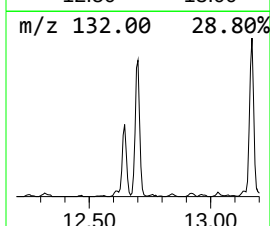
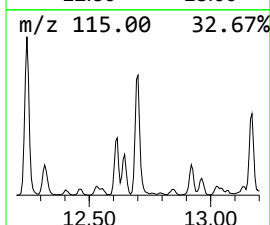
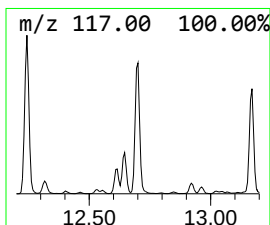
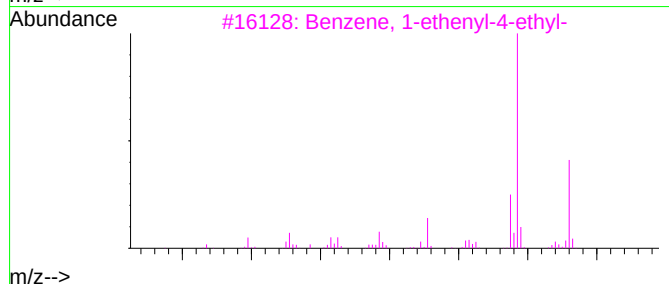
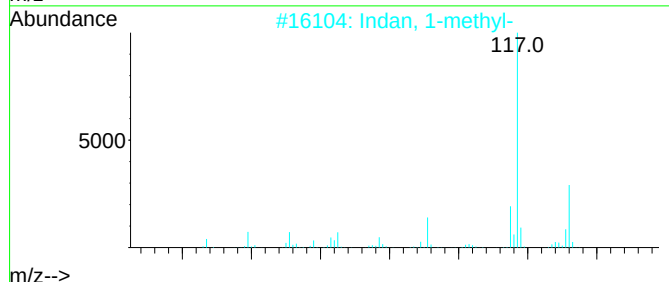
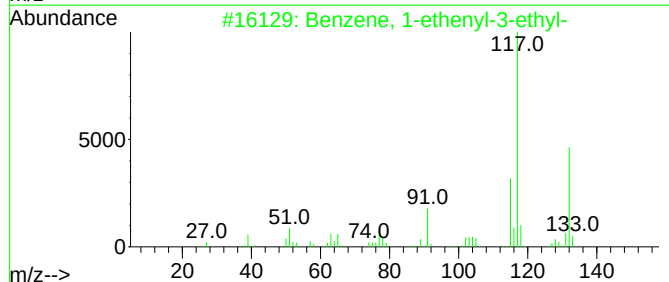
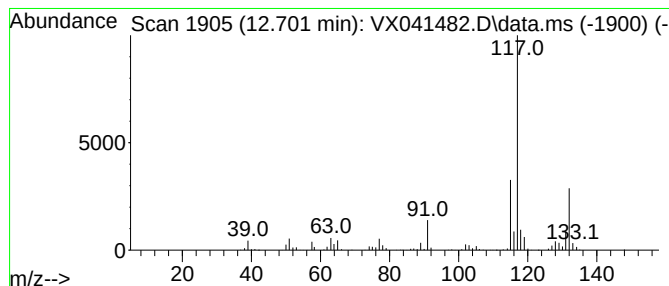
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	21.67 ug/l	339112	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
2	Indan, 1-methyl-	132	C10H12	000767-58-8	87
3	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
4	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	86
5	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	86



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051324\
Data File : VX041482.D
Acq On : 13 May 2024 11:13
Operator : JC/MD
Sample : P2413-01ME
Misc : 4.99g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 7 Sample Multiplier: 1

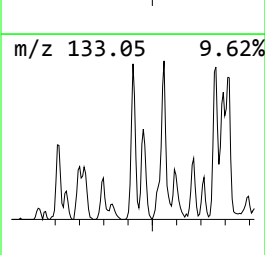
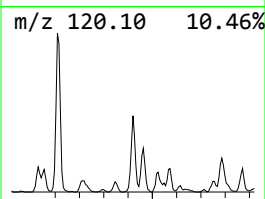
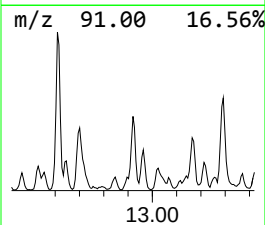
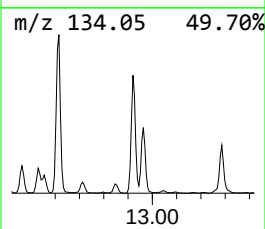
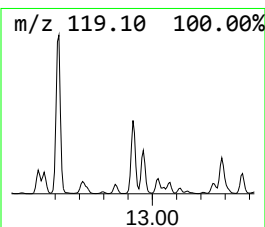
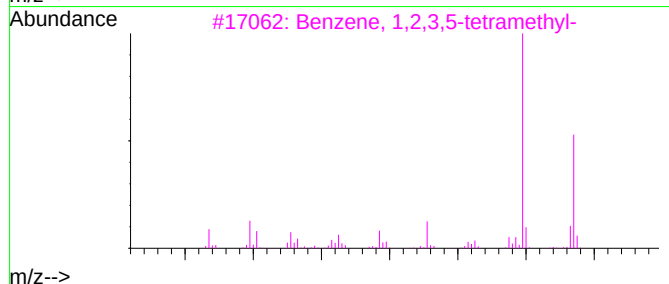
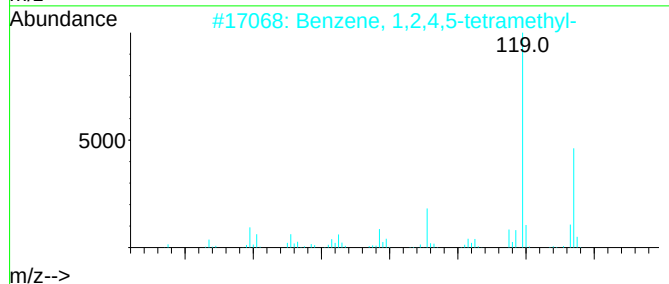
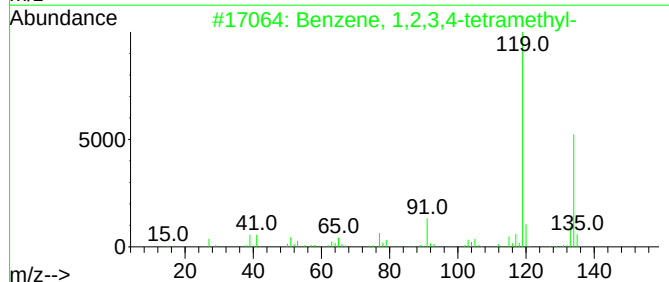
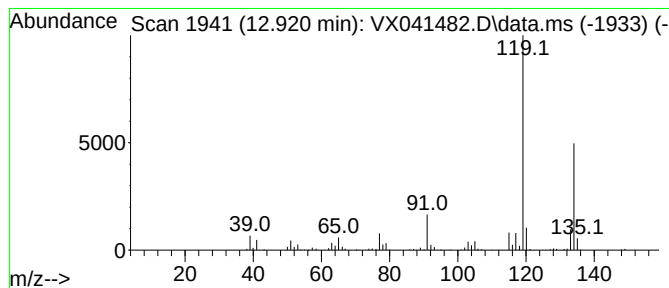
Instrument :
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ClientSampleId :
WC-CCME

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X043024W.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.920	17.75 ug/l	277728	1,4-Dichlorobenzene-d4	12.024	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051324\
Data File : VX041482.D
Acq On : 13 May 2024 11:13
Operator : JC/MD
Sample : P2413-01ME
Misc : 4.99g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WC-CCME

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X043024W.M
Quant Title : SW846 8260

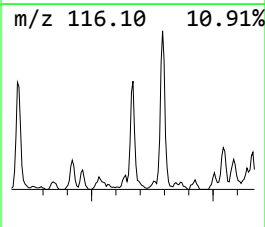
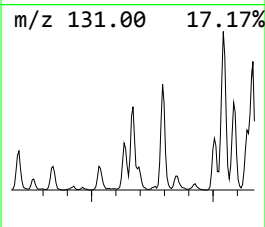
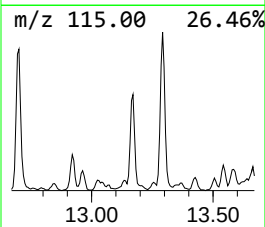
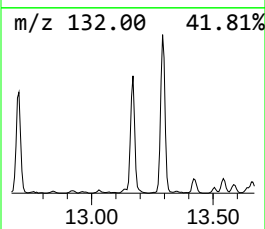
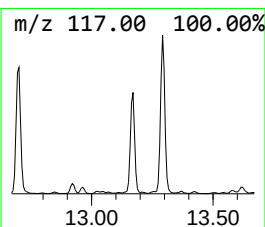
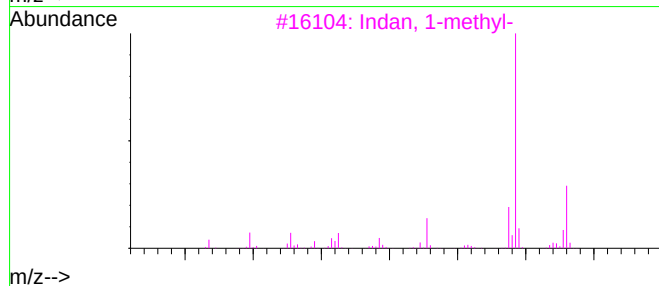
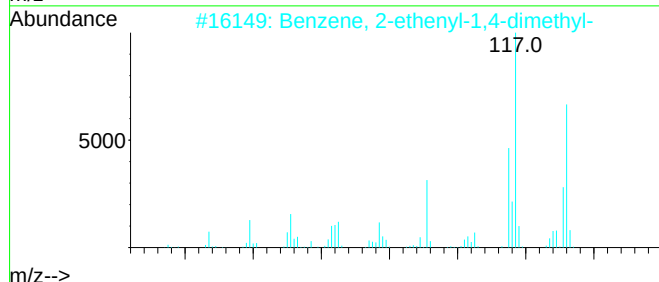
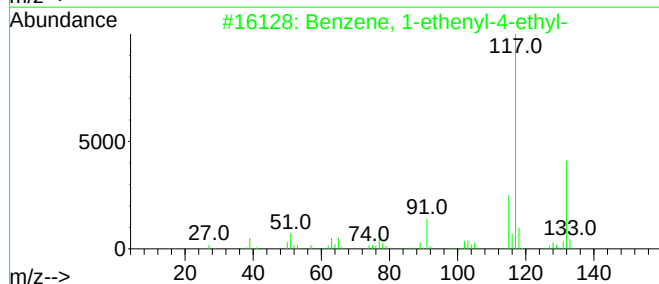
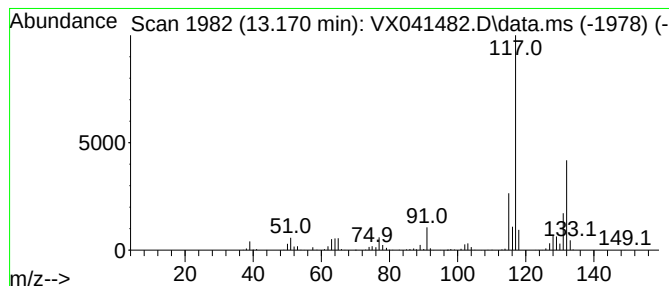
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.170	16.56 ug/l	259107	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
3	Indan, 1-methyl-	132	C10H12	000767-58-8	91
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051324\
Data File : VX041482.D
Acq On : 13 May 2024 11:13
Operator : JC/MD
Sample : P2413-01ME
Misc : 4.99g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 7 Sample Multiplier: 1

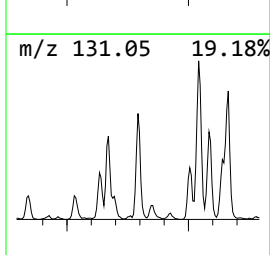
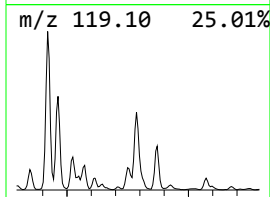
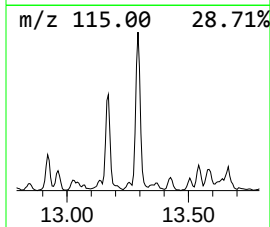
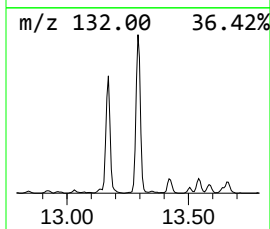
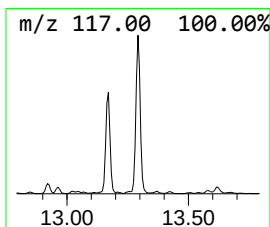
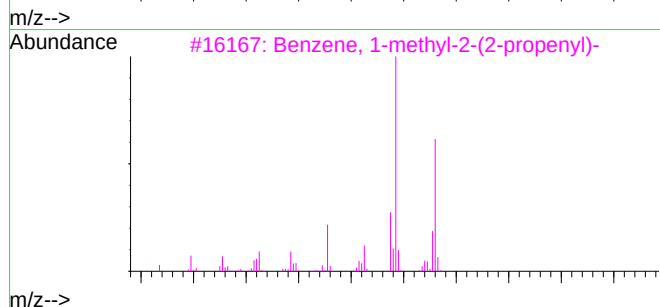
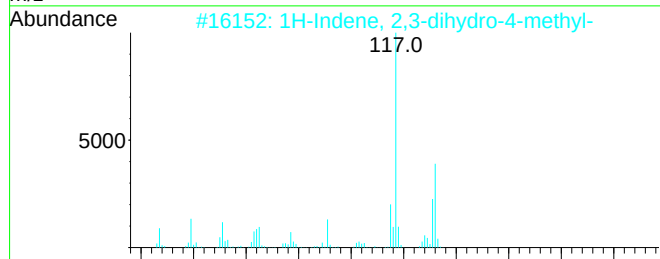
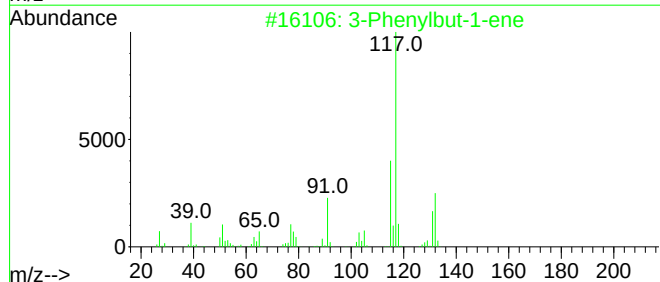
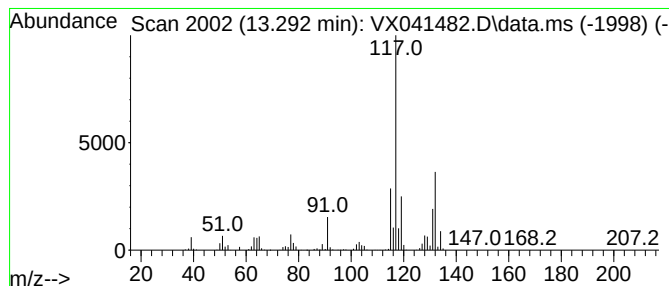
Instrument :
MSVOA_X
ClientSampleId :
WC-CCME

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X043024W.M
Quant Title : SW846 8260

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

Peak Number 10 3-Phenylbut-1-ene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.292	31.85 ug/l	498381	1,4-Dichlorobenzene-d4		12.024	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	3-Phenylbut-1-ene		132	C10H12	000934-10-1	87
2	1H-Indene, 2,3-dihydro-4-methyl-		132	C10H12	000824-22-6	81
3	Benzene, 1-methyl-2-(2-propenyl)-		132	C10H12	001587-04-8	74
4	Benzene, 1-ethenyl-4-ethyl-		132	C10H12	003454-07-7	70
5	Indan, 1-methyl-		132	C10H12	000767-58-8	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051324\
Data File : VX041482.D
Acq On : 13 May 2024 11:13
Operator : JC/MD
Sample : P2413-01ME
Misc : 4.99g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WC-CCME

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X043024W.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Hexane, 3-methyl-	5.806	25.0	ug/l	309786	1	5.544	619609	50.0
1-Butanol, 2-me...	6.239	12.4	ug/l	149520	2	6.757	602809	50.0
Heptane, 3-methyl-	8.433	15.1	ug/l	222041	3	10.055	737306	50.0
Octane, 2-methyl-	9.902	11.8	ug/l	173848	3	10.055	737306	50.0
Benzene, 1-prop...	12.244	33.9	ug/l	531198	4	12.024	782473	50.0
p-Cymene	12.609	32.1	ug/l	502636	4	12.024	782473	50.0
Benzene, 1-ethe...	12.701	21.7	ug/l	339112	4	12.024	782473	50.0
Benzene, 1,2,3,...	12.920	17.8	ug/l	277728	4	12.024	782473	50.0
Benzene, 1-ethe...	13.170	16.6	ug/l	259107	4	12.024	782473	50.0
3-Phenylbut-1-ene	13.292	31.9	ug/l	498381	4	12.024	782473	50.0