

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120224\
 Data File : VX044059.D
 Acq On : 02 Dec 2024 12:35
 Operator : JC/MD
 Sample : P5021-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-06

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112124W.M
 Title : SW846 8260

Signal : TIC: VX044059.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.587	69	82	84	rBV3	16870	59914	3.54%	0.316%
2	2.770	269	276	279	rBV3	13949	29780	1.76%	0.157%
3	2.812	279	283	288	rVB2	12397	24089	1.42%	0.127%
4	3.093	322	329	341	rVB2	22923	55717	3.29%	0.294%
5	3.721	423	432	441	rBV2	13928	33699	1.99%	0.178%
6	4.202	501	511	518	rBV4	18531	53083	3.13%	0.280%
7	4.288	518	525	537	rVB3	41908	114563	6.76%	0.605%
8	5.379	693	704	713	rBV	102366	306681	18.10%	1.619%
9	5.543	722	731	738	rBV	143725	416495	24.59%	2.199%
10	5.818	766	776	787	rBV2	19960	62007	3.66%	0.327%
11	5.952	788	798	806	rVV	118879	329466	19.45%	1.739%
12	6.037	806	812	820	rVV2	48424	134483	7.94%	0.710%
13	6.147	824	830	831	rVV2	14437	26303	1.55%	0.139%
14	6.238	835	845	856	rVV2	132562	436355	25.76%	2.304%
15	6.348	856	863	876	rVB5	14231	42415	2.50%	0.224%
16	6.757	921	930	940	rBV	260644	655138	38.67%	3.459%
17	7.165	987	997	1004	rVB3	13640	32754	1.93%	0.173%
18	7.360	1015	1029	1043	rVB5	38844	124034	7.32%	0.655%
19	7.488	1043	1050	1056	rBV3	15517	35515	2.10%	0.187%
20	7.555	1056	1061	1065	rVV3	15463	36196	2.14%	0.191%
21	7.598	1065	1068	1072	rVV2	12967	23461	1.38%	0.124%
22	7.653	1072	1077	1084	rVB3	10888	23167	1.37%	0.122%
23	7.903	1111	1118	1123	rVV4	22698	61390	3.62%	0.324%
24	7.976	1123	1130	1142	rVB	123238	256886	15.16%	1.356%
25	8.104	1142	1151	1160	rBV2	190379	433341	25.58%	2.288%
26	8.183	1160	1164	1174	rVB3	18674	40369	2.38%	0.213%
27	8.433	1200	1205	1212	rBV	74419	134995	7.97%	0.713%
28	8.500	1212	1216	1224	rVB4	26230	47930	2.83%	0.253%
29	8.592	1224	1231	1232	rBV3	21475	40084	2.37%	0.212%
30	8.647	1234	1240	1247	rVV	519884	840082	49.59%	4.435%
31	8.714	1248	1251	1258	rVB4	18513	31717	1.87%	0.167%
32	8.811	1263	1267	1277	rVB	14756	29776	1.76%	0.157%
33	9.098	1307	1314	1318	rVB4	12418	21140	1.25%	0.112%
34	9.159	1319	1324	1328	rBV	24640	37794	2.23%	0.200%
35	9.244	1333	1338	1344	rVB	24295	38306	2.26%	0.202%
36	9.451	1364	1372	1377	rVV5	32373	66226	3.91%	0.350%

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37	9.945	1448	1453	1460	rBV	73692	110074	6.50%	0.581%
38	10.006	1460	1463	1465	rBV	16768	22227	1.31%	0.117%
39	10.049	1465	1470	1476	rVV	493612	696821	41.13%	3.679%
40	10.213	1490	1497	1500	rBV2	72348	130564	7.71%	0.689%
41	10.244	1500	1502	1507	rVV	34129	40246	2.38%	0.212%
42	10.299	1507	1511	1518	rVB	49619	74854	4.42%	0.395%
43	10.384	1518	1525	1529	rBV2	59456	105965	6.26%	0.559%
44	10.640	1563	1567	1573	rVB2	28175	40391	2.38%	0.213%
45	10.793	1582	1592	1599	rBV5	10829	35408	2.09%	0.187%
46	10.896	1599	1609	1614	rVV2	37876	75270	4.44%	0.397%
47	10.957	1614	1619	1625	rVV	314299	423270	24.99%	2.235%
48	11.012	1625	1628	1634	rVB2	27378	35982	2.12%	0.190%
49	11.079	1634	1639	1649	rBV	438080	617811	36.47%	3.262%
50	11.158	1649	1652	1659	rVB4	17312	37439	2.21%	0.198%
51	11.299	1668	1675	1682	rBV	653946	897790	53.00%	4.740%
52	11.451	1695	1700	1703	rVV5	21102	38230	2.26%	0.202%
53	11.494	1703	1707	1712	rVB2	22296	33498	1.98%	0.177%
54	11.609	1721	1726	1731	rBV	65747	86196	5.09%	0.455%
55	11.713	1738	1743	1744	rBV3	13457	20437	1.21%	0.108%
56	11.750	1744	1749	1754	rVB2	33677	58824	3.47%	0.311%
57	11.865	1763	1768	1769	rBV	36648	45476	2.68%	0.240%
58	11.890	1769	1772	1781	rVB	50776	75749	4.47%	0.400%
59	12.018	1788	1793	1800	rVV	544668	688524	40.64%	3.635%
60	12.085	1800	1804	1808	rVB	43431	49515	2.92%	0.261%
61	12.176	1816	1819	1823	rVB	23200	26076	1.54%	0.138%
62	12.243	1823	1830	1836	rBV	1342680	1694002	100.00%	8.943%
63	12.311	1837	1841	1850	rVV2	162345	246840	14.57%	1.303%
64	12.402	1851	1856	1861	rVV2	128606	160399	9.47%	0.847%
65	12.469	1861	1867	1872	rVB3	52916	94518	5.58%	0.499%
66	12.530	1872	1877	1884	rBV6	17586	32627	1.93%	0.172%
67	12.609	1885	1890	1894	rBV	1096501	1301536	76.83%	6.871%
68	12.640	1894	1895	1899	rVV	55036	55297	3.26%	0.292%
69	12.695	1899	1904	1912	rVB2	326048	532767	31.45%	2.813%
70	12.835	1923	1927	1932	rVB3	20700	32460	1.92%	0.171%
71	12.920	1932	1941	1947	rBV	673477	838810	49.52%	4.428%
72	12.963	1947	1948	1953	rVV3	15101	19724	1.16%	0.104%
73	13.024	1953	1958	1966	rVV2	57801	93330	5.51%	0.493%
74	13.115	1966	1973	1975	rVV3	42902	79494	4.69%	0.420%
75	13.170	1978	1982	1988	rVV	207781	279982	16.53%	1.478%
76	13.231	1988	1992	1993	rVV	26861	29977	1.77%	0.158%

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77	13.255	1993	1996	1998	rVV	53848	75904	4.48%	0.401%
78	13.286	1998	2001	2007	rVV2	444073	660581	39.00%	3.487%
79	13.341	2007	2010	2012	rVV2	57214	73916	4.36%	0.390%
80	13.365	2012	2014	2020	rVV	52282	79624	4.70%	0.420%
81	13.420	2020	2023	2028	rVB	49588	62240	3.67%	0.329%
82	13.505	2033	2037	2039	rVV3	22826	32524	1.92%	0.172%
83	13.542	2039	2043	2046	rVV	93783	139491	8.23%	0.736%
84	13.579	2046	2049	2053	rVV2	133563	178956	10.56%	0.945%
85	13.640	2053	2059	2060	rVV2	48440	80036	4.72%	0.423%
86	13.670	2060	2064	2070	rVV2	78027	146415	8.64%	0.773%
87	13.774	2074	2081	2091	rVV	999950	1244532	73.47%	6.570%
88	13.926	2099	2106	2110	rBV3	32154	55915	3.30%	0.295%
89	13.969	2110	2113	2117	rVB2	24649	31096	1.84%	0.164%
90	14.072	2125	2130	2136	rBV	74232	92848	5.48%	0.490%
91	14.213	2148	2153	2159	rBV2	58695	94570	5.58%	0.499%
92	14.316	2166	2170	2175	rBV	99316	120275	7.10%	0.635%
93	14.371	2175	2179	2183	rVV	45250	57581	3.40%	0.304%
94	14.597	2212	2216	2219	rBV2	16788	27140	1.60%	0.143%
95	14.633	2219	2222	2227	rVB	96334	120828	7.13%	0.638%
96	14.780	2240	2246	2256	rBV	202497	273513	16.15%	1.444%
97	15.481	2355	2361	2370	rVB4	9253	20636	1.22%	0.109%
98	15.615	2376	2383	2386	rBV2	24406	39350	2.32%	0.208%
99	16.322	2491	2499	2508	rBV	67281	119286	7.04%	0.630%
100	16.688	2549	2559	2565	rBV	9474	21044	1.24%	0.111%

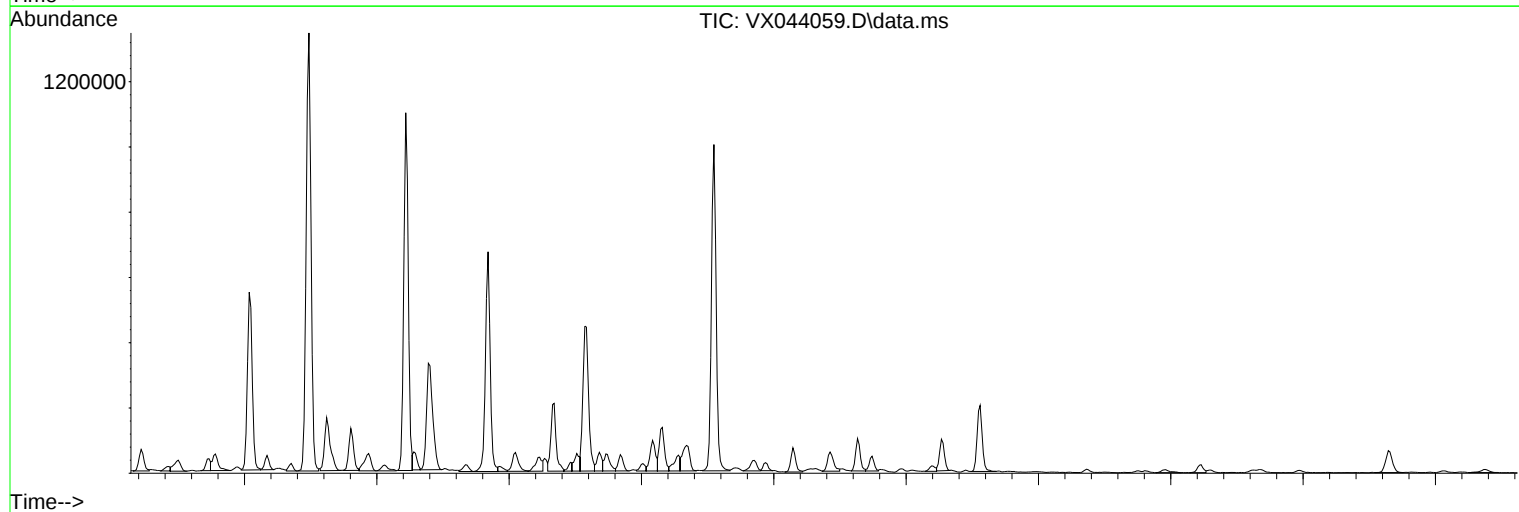
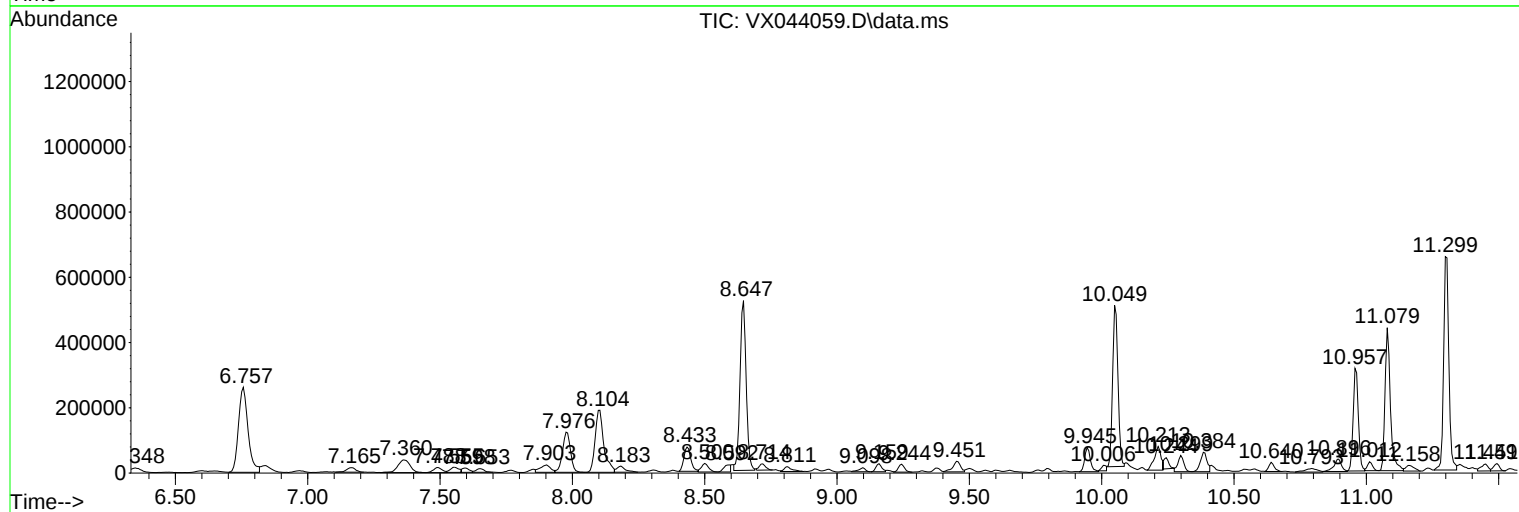
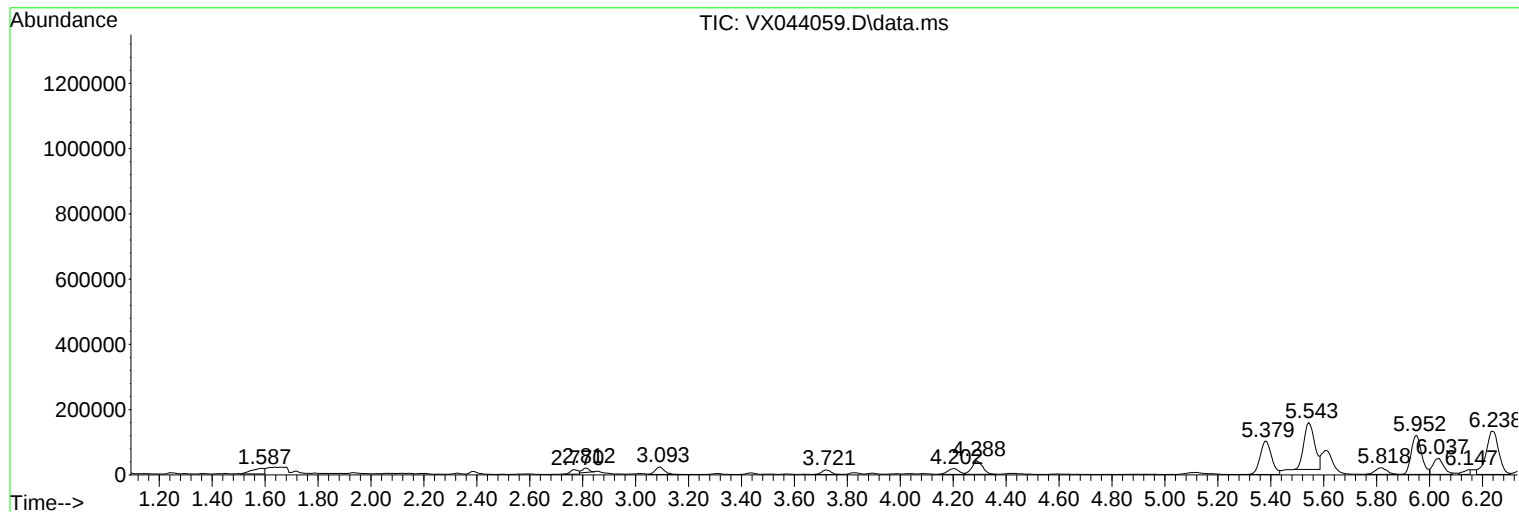
Sum of corrected areas: 18942052

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Instrument :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112124W.M
Quant Title : SW846 8260

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



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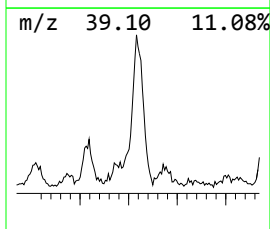
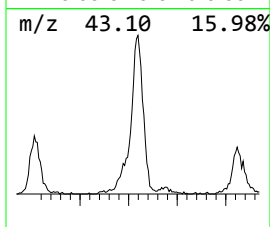
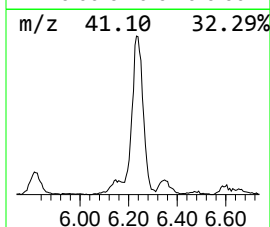
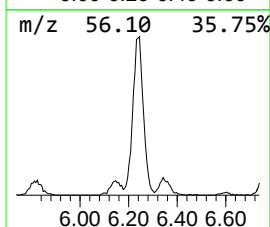
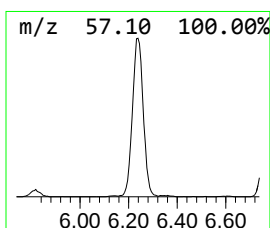
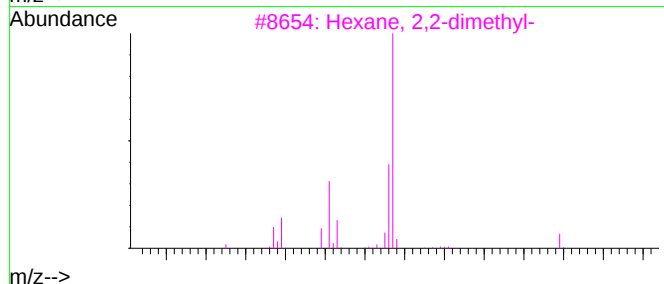
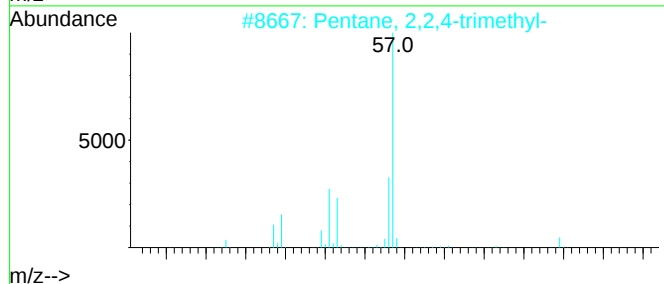
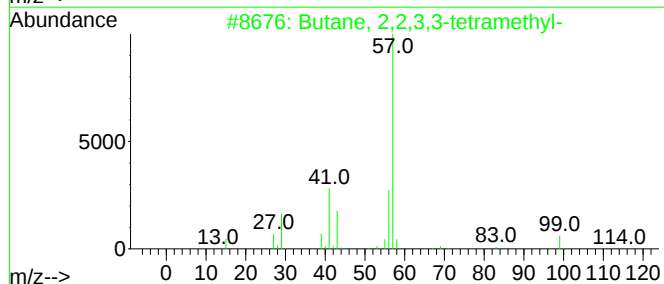
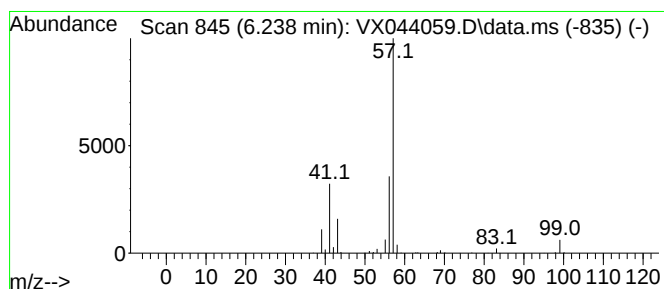
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2,2,3,3-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.238	33.30 ug/l	436355	1,4-Difluorobenzene	6.757

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	78
2	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
3	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	72
4	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	64
5	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	45



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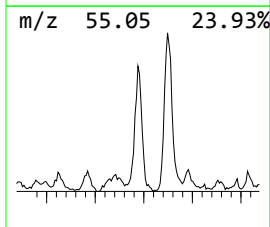
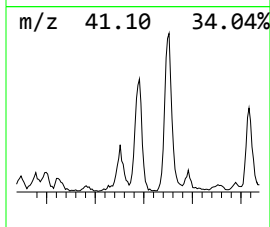
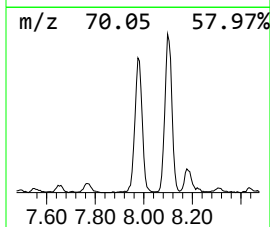
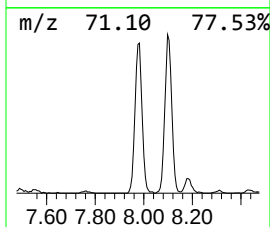
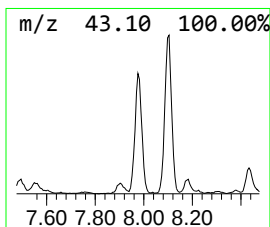
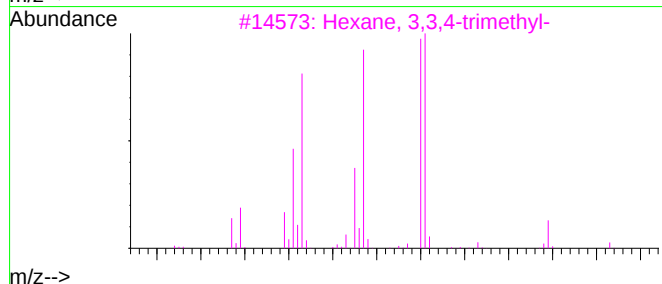
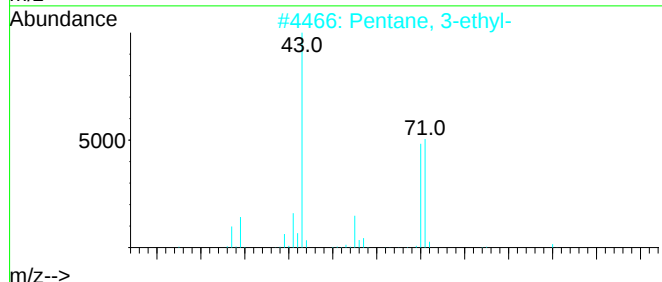
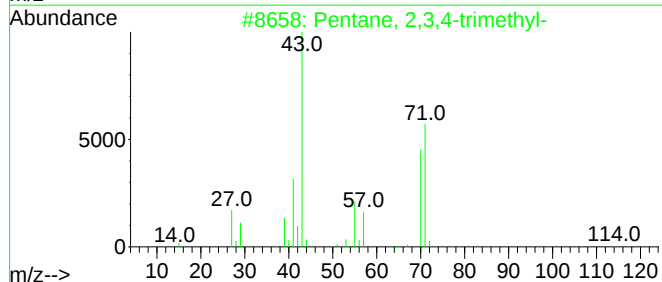
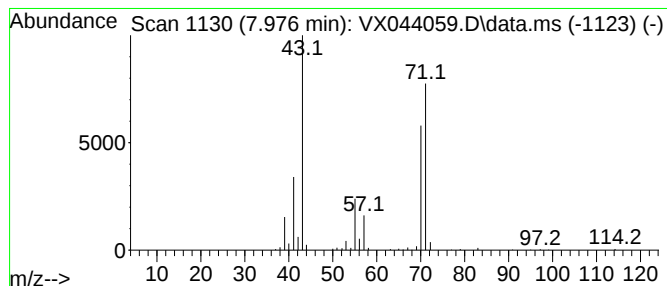
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 2 Pentane, 2,3,4-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.976	19.61 ug/l	256886	1,4-Difluorobenzene	6.757

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	91
2	Pentane, 3-ethyl-	100	C7H16	000617-78-7	83
3	Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	78
4	Heptane, 4-methyl-	114	C8H18	000589-53-7	78
5	Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	56



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Quant Title : SW846 8260

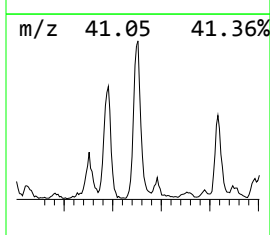
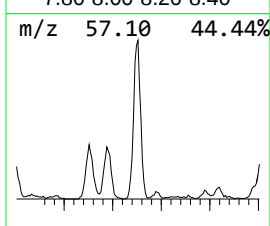
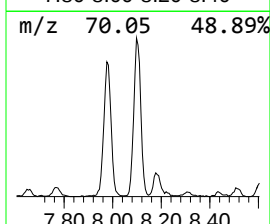
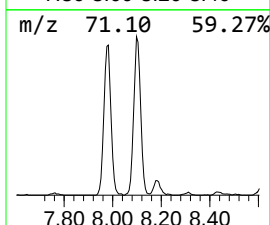
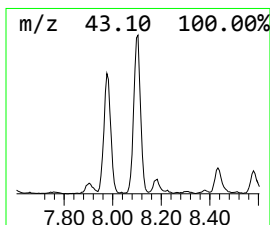
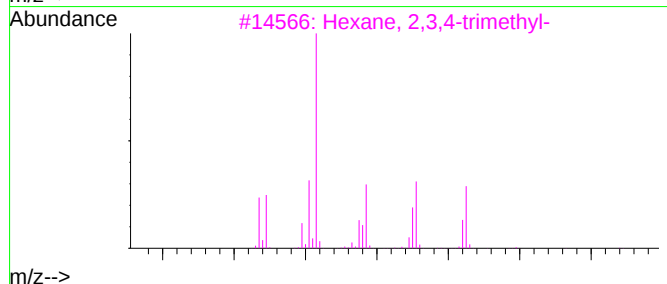
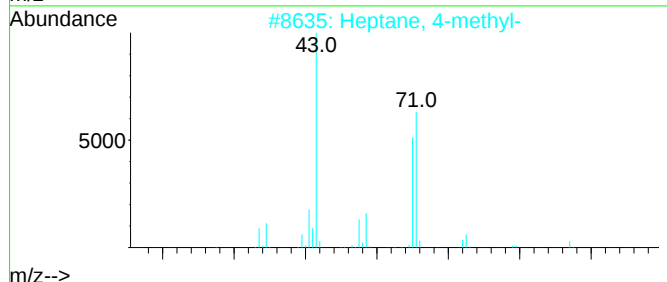
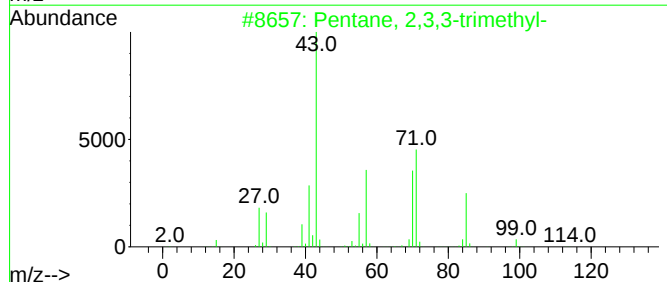
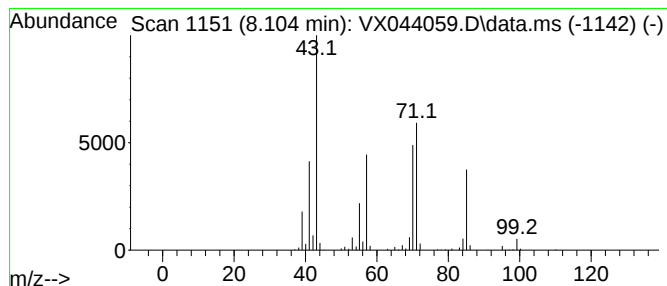
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 3 Pentane, 2,3,3-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.104	33.07 ug/l	433341	1,4-Difluorobenzene	6.757

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2	Heptane, 4-methyl-	114	C8H18	000589-53-7	64
3	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	59
4	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	59
5	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120224\
Data File : VX044059.D
Acq On : 02 Dec 2024 12:35
Operator : JC/MD
Sample : P5021-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-06

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112124W.M
Quant Title : SW846 8260

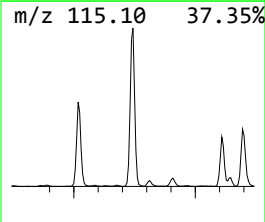
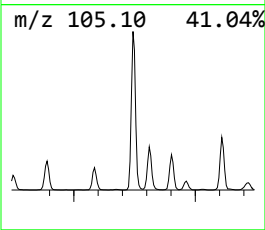
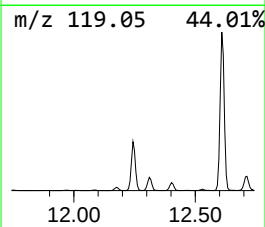
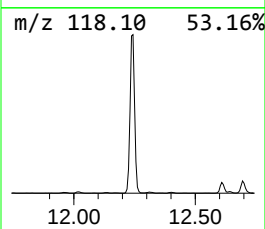
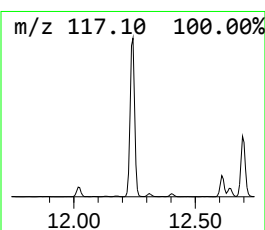
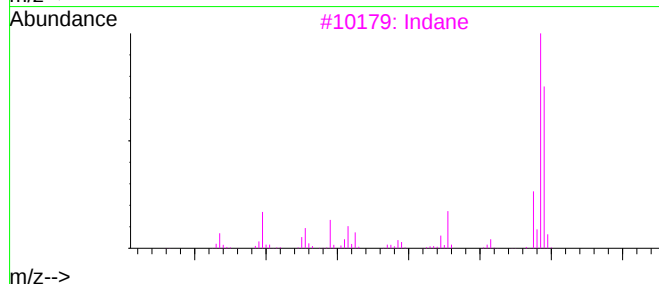
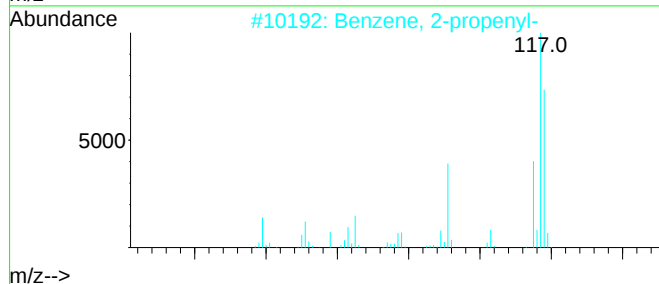
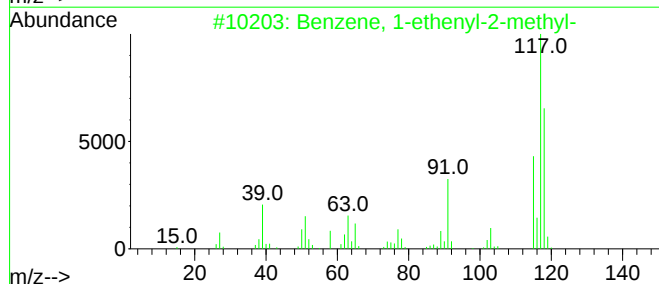
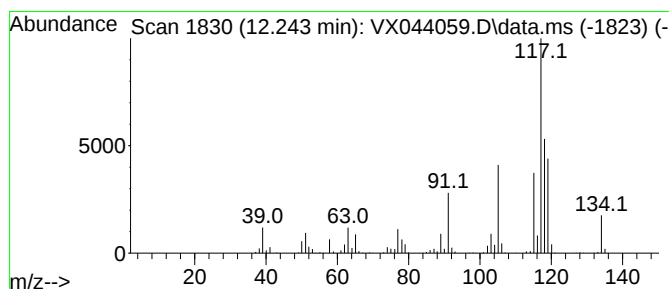
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TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1-ethenyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.243	123.02 ug/l	1694000	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	80
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	60
3	Indane	118	C9H10	000496-11-7	60
4	Benzene, 1-propenyl-	118	C9H10	000637-50-3	60
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120224\
Data File : VX044059.D
Acq On : 02 Dec 2024 12:35
Operator : JC/MD
Sample : P5021-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-06

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112124W.M
Quant Title : SW846 8260

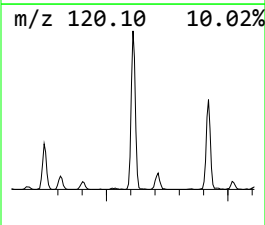
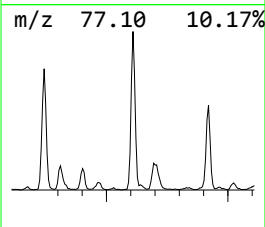
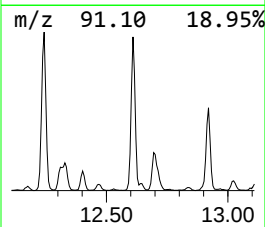
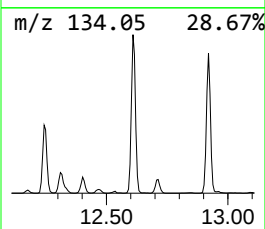
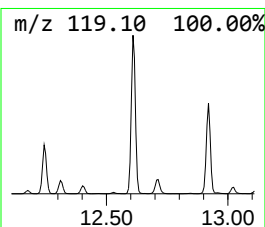
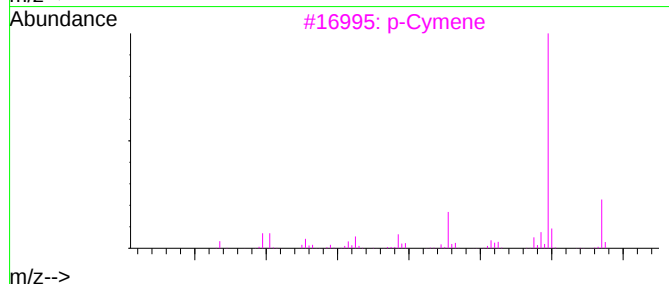
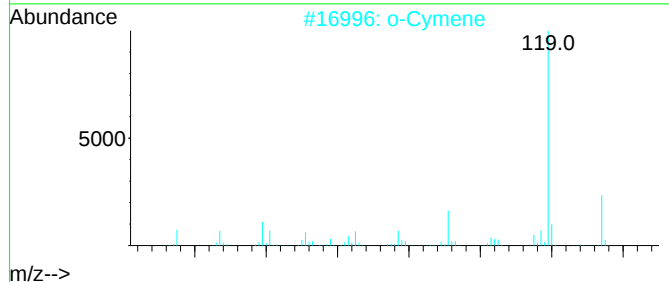
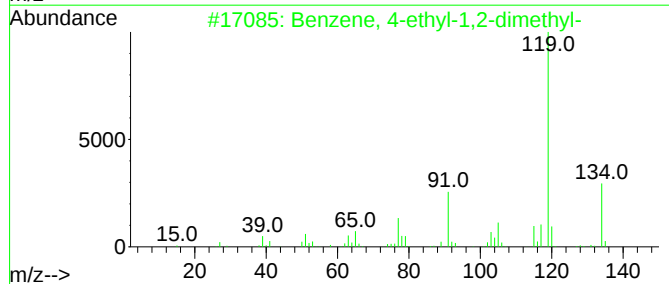
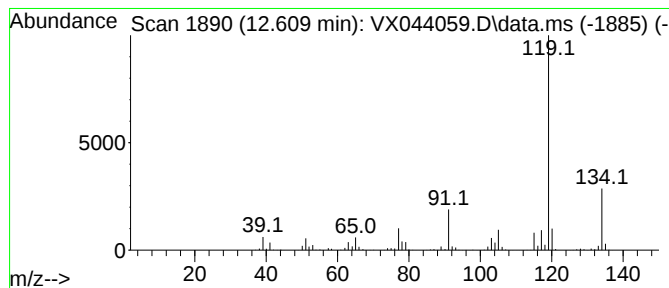
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.609	94.52 ug/l	1301540	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2	o-Cymene	134	C10H14	000527-84-4	95
3	p-Cymene	134	C10H14	000099-87-6	95
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120224\
Data File : VX044059.D
Acq On : 02 Dec 2024 12:35
Operator : JC/MD
Sample : P5021-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-06

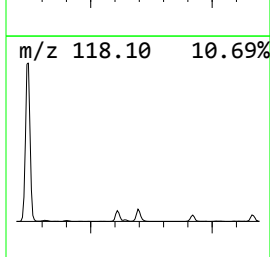
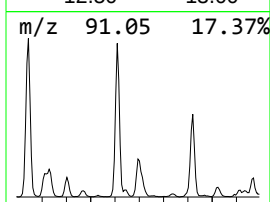
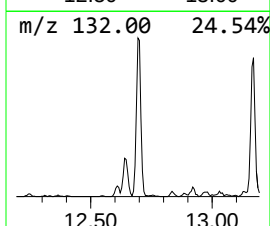
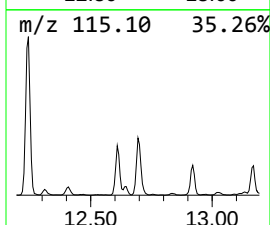
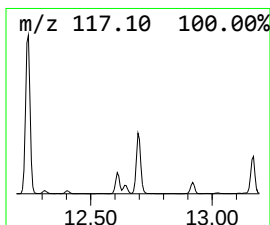
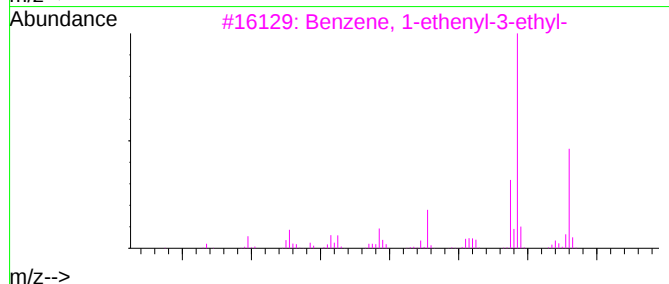
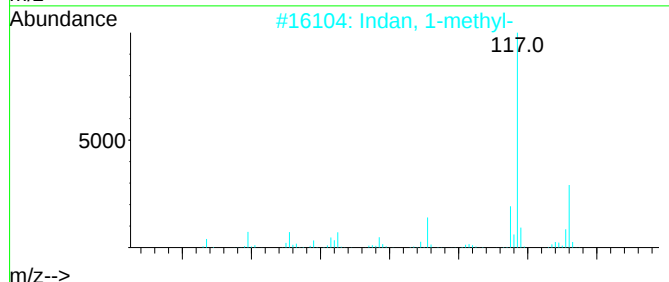
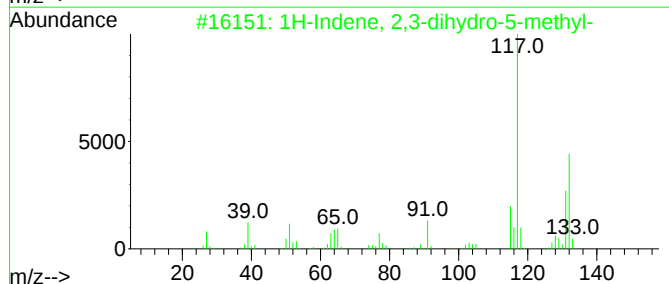
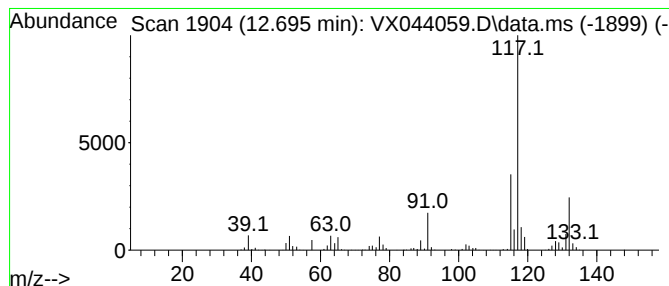
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 6 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.695	38.69 ug/l	532767	1,4-Dichlorobenzene-d4	12.024	
Hit# of 5	Tentative ID		MW MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-		132 C10H12	000874-35-1	87
2	Indan, 1-methyl-		132 C10H12	000767-58-8	87
3	Benzene, 1-ethenyl-3-ethyl-		132 C10H12	007525-62-4	87
4	(E)-1-Phenyl-1-butene		132 C10H12	001005-64-7	87
5	Benzene, (2-methyl-2-propenyl)-		132 C10H12	003290-53-7	86



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120224\
Data File : VX044059.D
Acq On : 02 Dec 2024 12:35
Operator : JC/MD
Sample : P5021-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-06

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112124W.M
Quant Title : SW846 8260

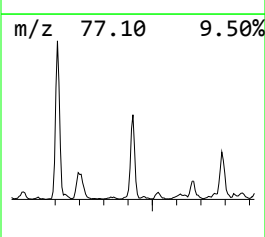
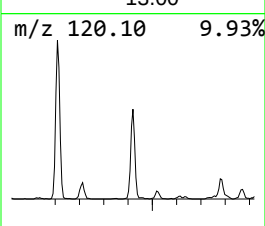
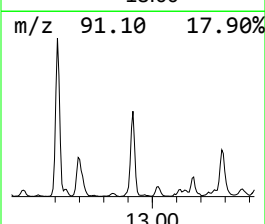
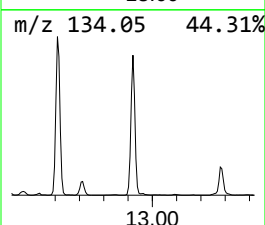
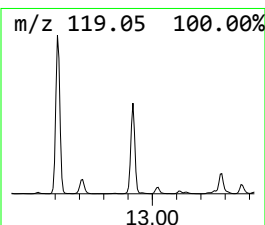
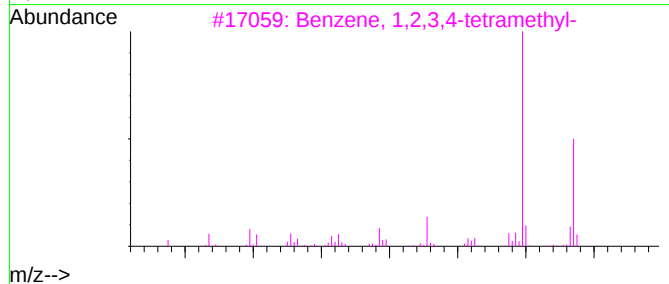
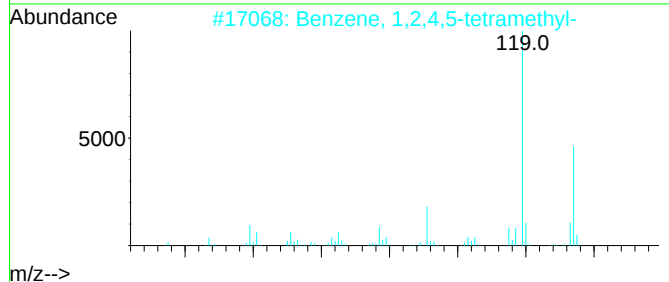
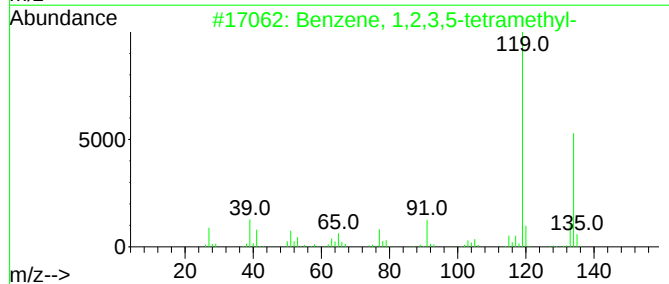
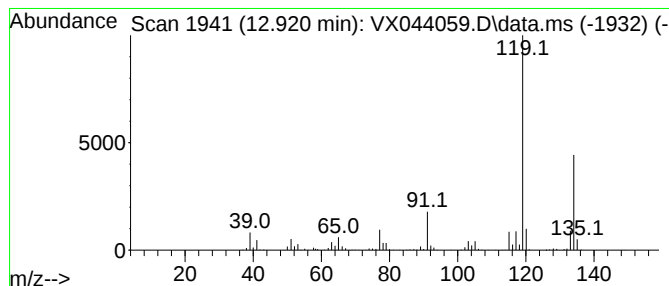
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.920	60.91 ug/l	838810	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120224\
Data File : VX044059.D
Acq On : 02 Dec 2024 12:35
Operator : JC/MD
Sample : P5021-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-06

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112124W.M
Quant Title : SW846 8260

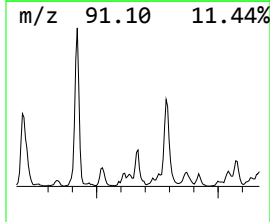
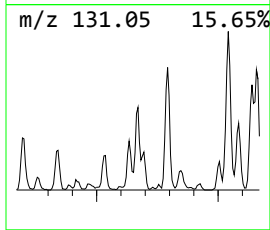
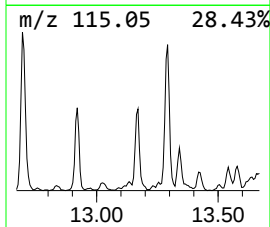
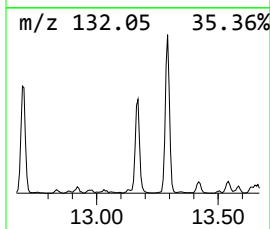
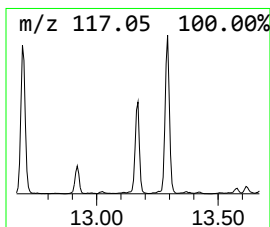
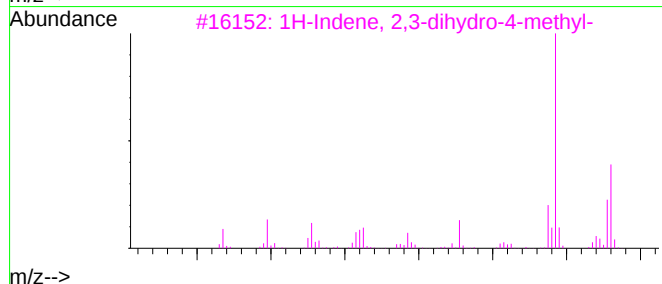
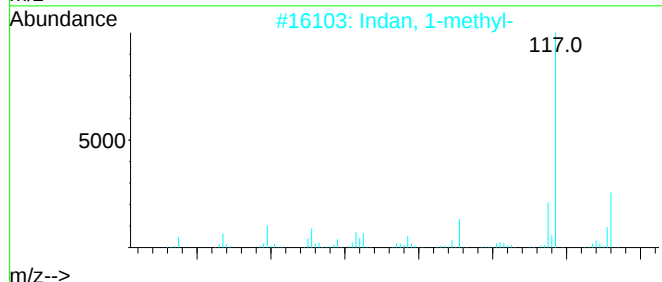
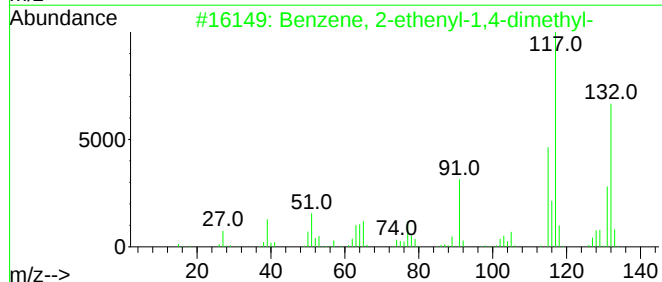
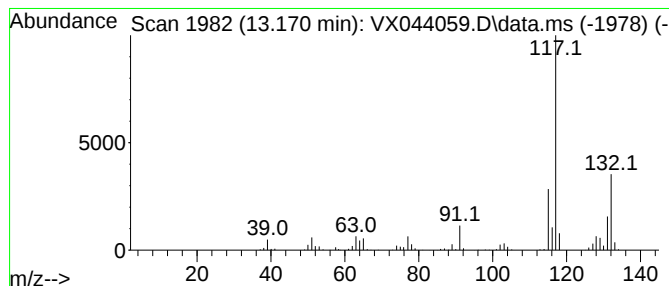
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.170	20.33 ug/l	279982	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
2	Indan, 1-methyl-	132	C10H12	000767-58-8	90
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
4	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120224\
Data File : VX044059.D
Acq On : 02 Dec 2024 12:35
Operator : JC/MD
Sample : P5021-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-06

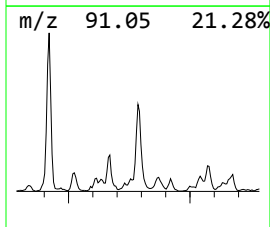
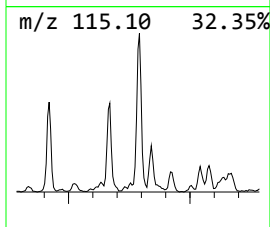
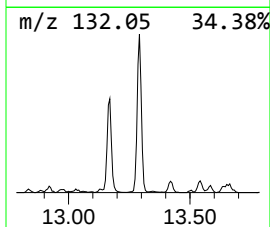
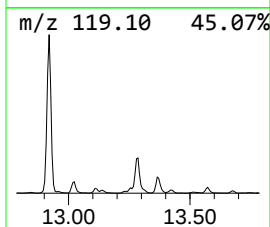
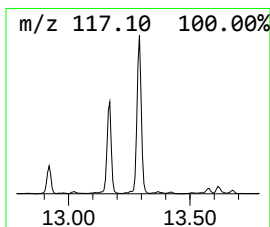
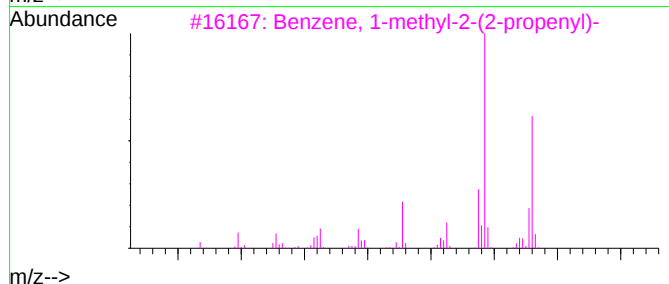
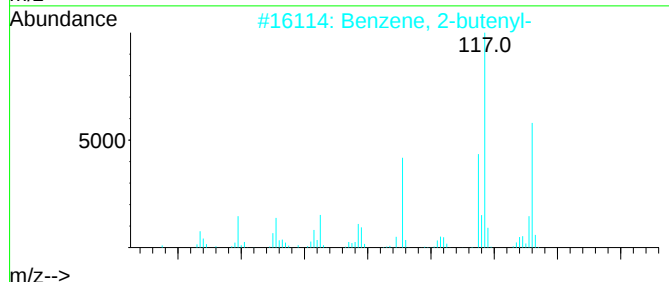
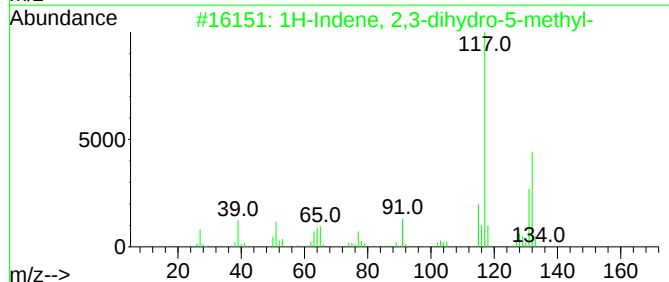
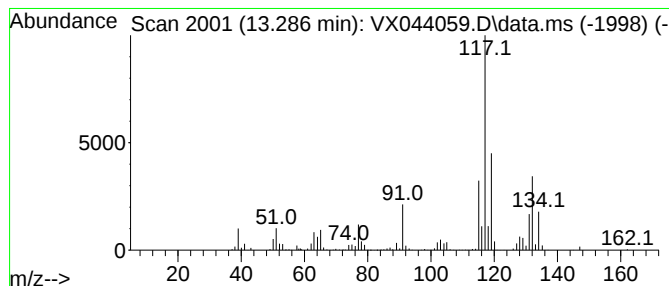
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, 2-butenyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.286	47.97 ug/l	660581	1,4-Dichlorobenzene-d4		12.024	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-		132	C10H12	000874-35-1	70
2	Benzene, 2-butenyl-		132	C10H12	001560-06-1	70
3	Benzene, 1-methyl-2-(2-propenyl)-		132	C10H12	001587-04-8	70
4	Benzene, 2-ethenyl-1,4-dimethyl-		132	C10H12	002039-89-6	70
5	Benzene, 1-ethenyl-4-ethyl-		132	C10H12	003454-07-7	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120224\
Data File : VX044059.D
Acq On : 02 Dec 2024 12:35
Operator : JC/MD
Sample : P5021-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-06

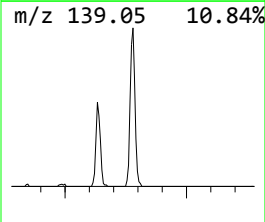
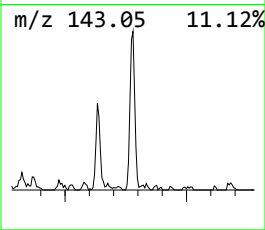
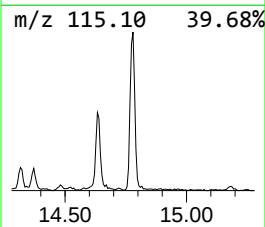
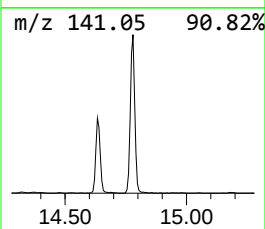
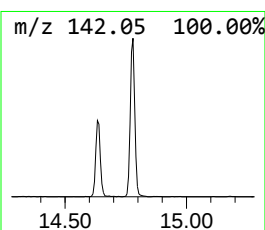
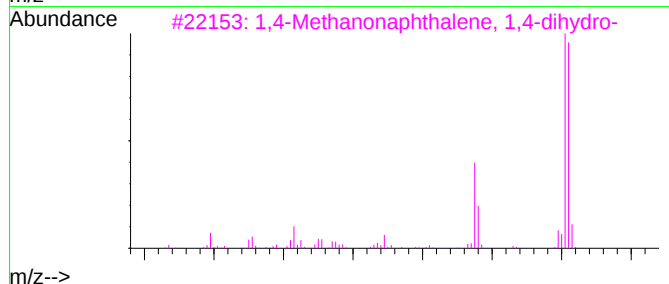
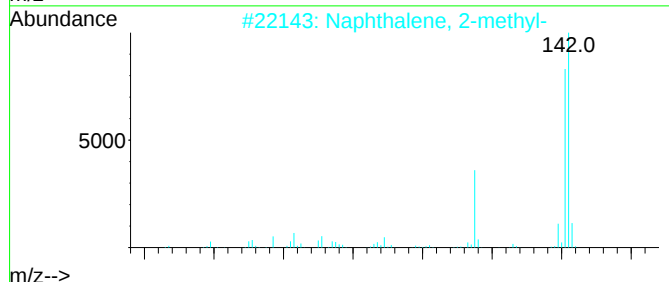
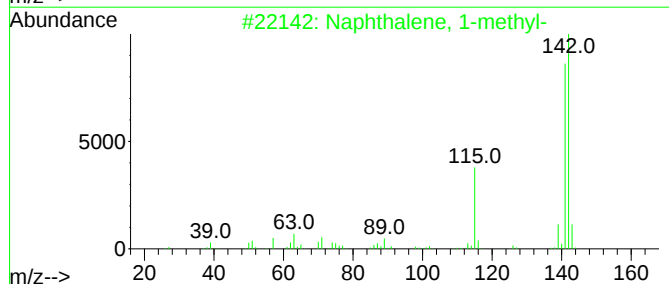
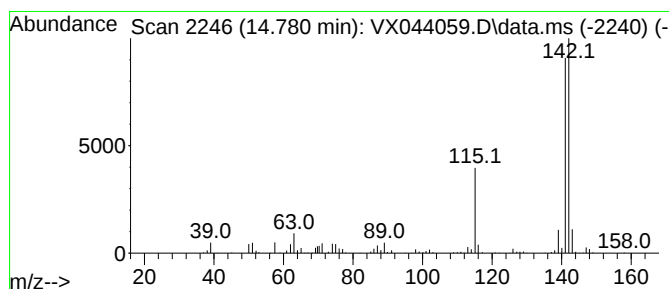
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.780	19.86 ug/l	273513	1,4-Dichlorobenzene-d4		12.024	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Naphthalene, 1-methyl-		142	C11H10	000090-12-0	96
2	Naphthalene, 2-methyl-		142	C11H10	000091-57-6	96
3	1,4-Methanonaphthalene, 1,4-dihy...		142	C11H10	004453-90-1	94
4	Benzocycloheptatriene		142	C11H10	000264-09-5	91
5	1H-Indene, 1-ethylidene-		142	C11H10	002471-83-2	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120224\
Data File : VX044059.D
Acq On : 02 Dec 2024 12:35
Operator : JC/MD
Sample : P5021-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-06

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112124W.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
Butane, 2,2,3,3...	6.238	33.3	ug/l	436355	2	6.757	655138	50.0
Pentane, 2,3,4-...	7.976	19.6	ug/l	256886	2	6.757	655138	50.0
Pentane, 2,3,3-...	8.104	33.1	ug/l	433341	2	6.757	655138	50.0
Benzene, 1-ethe...	12.243	123.0	ug/l	1694000	4	12.024	688524	50.0
Benzene, 4-ethy...	12.609	94.5	ug/l	1301540	4	12.024	688524	50.0
1H-Indene, 2,3-...	12.695	38.7	ug/l	532767	4	12.024	688524	50.0
Benzene, 1,2,3,...	12.920	60.9	ug/l	838810	4	12.024	688524	50.0
Benzene, 2-ethe...	13.170	20.3	ug/l	279982	4	12.024	688524	50.0
Benzene, 2-bute...	13.286	48.0	ug/l	660581	4	12.024	688524	50.0
Naphthalene, 1-...	14.780	19.9	ug/l	273513	4	12.024	688524	50.0