

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060425\
 Data File : VX046495.D
 Acq On : 04 Jun 2025 13:05
 Operator : JC/MD
 Sample : Q2169-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 303-PPR-1

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\82X050525W.M
 Title : SW846 8260

Signal : TIC: VX046495.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.380	207	212	225	rVB	25199	41677	1.14%	0.119%
2	5.385	694	705	714	rBV	59654	175862	4.81%	0.502%
3	5.544	723	731	751	rVB	85273	248918	6.81%	0.711%
4	5.952	788	798	806	rBV	67879	180803	4.95%	0.516%
5	6.037	806	812	825	rVB2	25371	69576	1.90%	0.199%
6	6.757	922	930	948	rBV	149659	367768	10.07%	1.050%
7	7.373	1020	1031	1042	rVB2	24836	61548	1.68%	0.176%
8	8.574	1222	1228	1234	rBV	101946	172760	4.73%	0.493%
9	8.647	1234	1240	1246	rVV	331275	518574	14.19%	1.480%
10	8.714	1246	1251	1264	rVB	1028203	1602102	43.85%	4.574%
11	10.049	1465	1470	1480	rBV	328065	466130	12.76%	1.331%
12	10.189	1488	1493	1503	rBV	795538	1069044	29.26%	3.052%
13	10.299	1503	1511	1519	rVV	2733814	3653598	100.00%	10.431%
14	10.640	1561	1567	1581	rVB	1657453	2101463	57.52%	6.000%
15	10.957	1614	1619	1627	rVB	104677	138508	3.79%	0.395%
16	11.079	1634	1639	1645	rBV	282109	356814	9.77%	1.019%
17	11.305	1670	1676	1682	rBV	227848	288037	7.88%	0.822%
18	11.378	1682	1688	1695	rVV	1084289	1750724	47.92%	4.998%
19	11.451	1695	1700	1715	rVB	509196	665579	18.22%	1.900%
20	11.567	1715	1719	1722	rBV	57200	71075	1.95%	0.203%
21	11.610	1722	1726	1739	rVB2	543032	835585	22.87%	2.386%
22	11.750	1744	1749	1754	rVV	2152729	2584208	70.73%	7.378%
23	11.823	1758	1761	1765	rVV	68306	98012	2.68%	0.280%
24	11.890	1769	1772	1776	rVB	56610	72108	1.97%	0.206%
25	11.969	1781	1785	1788	rVV	110595	139554	3.82%	0.398%
26	12.018	1788	1793	1799	rVV2	371577	548979	15.03%	1.567%
27	12.085	1799	1804	1814	rVV	968298	1174939	32.16%	3.354%
28	12.238	1822	1829	1833	rBV	710550	1084619	29.69%	3.097%
29	12.268	1833	1834	1837	rVV	212580	143464	3.93%	0.410%
30	12.317	1837	1842	1852	rVB3	301011	466972	12.78%	1.333%
31	12.408	1852	1857	1862	rBV3	45803	75076	2.05%	0.214%
32	12.457	1862	1865	1871	rVB	123653	162105	4.44%	0.463%
33	12.530	1871	1877	1879	rBV	207707	283413	7.76%	0.809%
34	12.555	1879	1881	1882	rVV	179345	170904	4.68%	0.488%
35	12.585	1882	1886	1889	rVV	2175391	2603926	71.27%	7.434%
36	12.640	1893	1895	1900	rVV	109989	162121	4.44%	0.463%

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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\82X050525W.M
 Title : SW846 8260

37	12.695	1900	1904	1912	rVB2	297399	449527	12.30%	1.283%
38	12.768	1912	1916	1918	rBV2	35461	47110	1.29%	0.134%
39	12.798	1918	1921	1925	rVV2	62763	86093	2.36%	0.246%
40	12.847	1925	1929	1935	rVB2	117574	155865	4.27%	0.445%
41	12.920	1935	1941	1944	rBV	175646	224506	6.14%	0.641%
42	12.963	1944	1948	1955	rVB	296532	366848	10.04%	1.047%
43	13.134	1973	1976	1978	rBV2	28377	36611	1.00%	0.105%
44	13.164	1978	1981	1985	rVB	282442	330820	9.05%	0.944%
45	13.250	1992	1995	1997	rBV2	39464	55491	1.52%	0.158%
46	13.286	1997	2001	2006	rVV2	682297	937265	25.65%	2.676%
47	13.335	2006	2009	2013	rVB4	52574	62293	1.70%	0.178%
48	13.420	2018	2023	2031	rVB	590385	722406	19.77%	2.062%
49	13.506	2032	2037	2039	rBV3	59392	78969	2.16%	0.225%
50	13.542	2039	2043	2046	rVV	76862	94548	2.59%	0.270%
51	13.585	2046	2050	2055	rVV2	105041	175332	4.80%	0.501%
52	13.658	2056	2062	2068	rVB2	111934	234745	6.43%	0.670%
53	13.725	2068	2073	2076	rBV2	49267	75025	2.05%	0.214%
54	13.774	2076	2081	2087	rVB	921153	1128341	30.88%	3.221%
55	13.853	2089	2094	2099	rVB	111115	159948	4.38%	0.457%
56	13.920	2099	2105	2110	rBV2	133859	191131	5.23%	0.546%
57	13.969	2110	2113	2117	rBV	50728	66350	1.82%	0.189%
58	14.073	2126	2130	2134	rBV	93695	116533	3.19%	0.333%
59	14.195	2145	2150	2151	rBV	195180	234420	6.42%	0.669%
60	14.225	2152	2155	2163	rVV	330117	479129	13.11%	1.368%
61	14.316	2167	2170	2175	rVB2	46040	52767	1.44%	0.151%
62	14.371	2175	2179	2184	rBV	98752	133251	3.65%	0.380%
63	14.475	2192	2196	2202	rBV2	213660	276407	7.57%	0.789%
64	14.633	2214	2222	2226	rBV	820257	1081591	29.60%	3.088%
65	14.670	2226	2228	2233	rVB2	64876	74409	2.04%	0.212%
66	14.774	2240	2245	2250	rVV	550435	721209	19.74%	2.059%
67	14.841	2254	2256	2261	rVB3	29322	36586	1.00%	0.104%
68	14.902	2261	2266	2272	rBV3	28331	56934	1.56%	0.163%
69	15.182	2307	2312	2313	rBV	63285	83688	2.29%	0.239%
70	15.200	2313	2315	2320	rVV	65792	91594	2.51%	0.261%
71	15.371	2338	2343	2346	rBV	63014	92001	2.52%	0.263%
72	15.475	2354	2360	2365	rBV	158116	249667	6.83%	0.713%
73	15.609	2375	2382	2385	rBV	212181	324792	8.89%	0.927%
74	15.645	2385	2388	2396	rVB	117260	179297	4.91%	0.512%
75	15.834	2409	2419	2434	rVB2	60758	162025	4.43%	0.463%
76	15.987	2437	2444	2454	rVB	35902	71228	1.95%	0.203%

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Instrument :
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ClientSampleId :
303-PPR-1

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

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Peak separation: 5

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Title : SW846 8260

77	16.084	2455	2460	2469	rBV3	21870	39695	1.09%	0.113%
78	16.322	2490	2499	2511	rBV3	26067	72477	1.98%	0.207%
79	16.475	2518	2524	2529	rBV	30305	56992	1.56%	0.163%
80	16.651	2548	2553	2570	rVB4	15806	54515	1.49%	0.156%

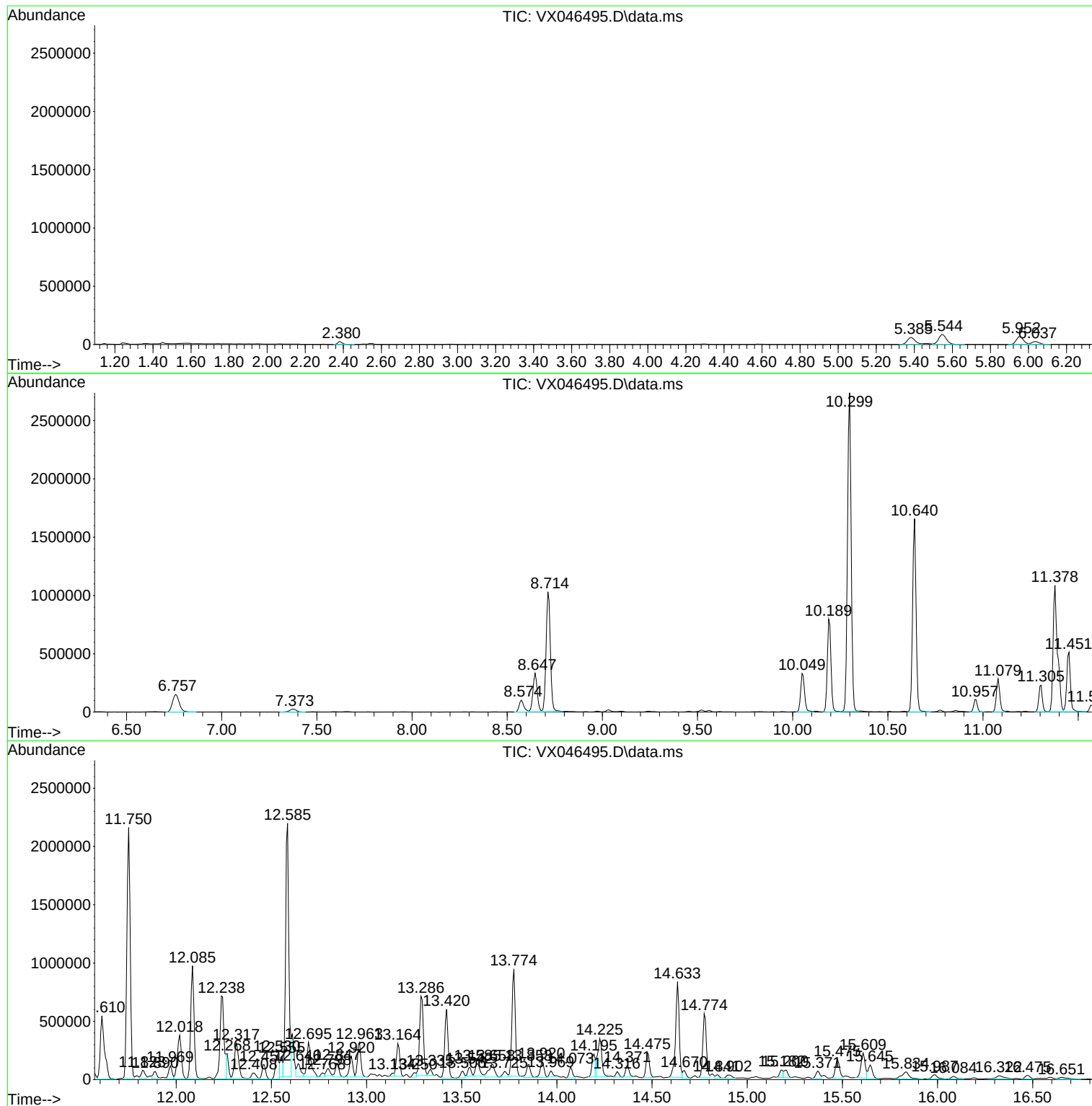
Sum of corrected areas: 35026976

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060425\
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Instrument :
MSVOA_X
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303-PPR-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X050525W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060425\
Data File : VX046495.D
Acq On : 04 Jun 2025 13:05
Operator : JC/MD
Sample : Q2169-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
303-PPR-1

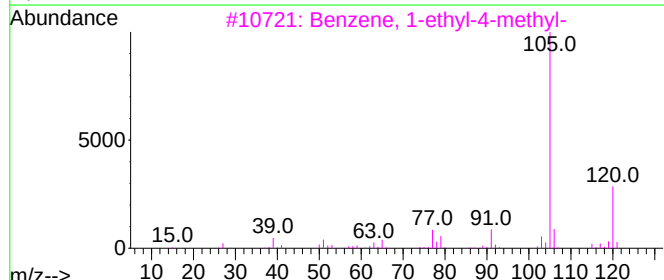
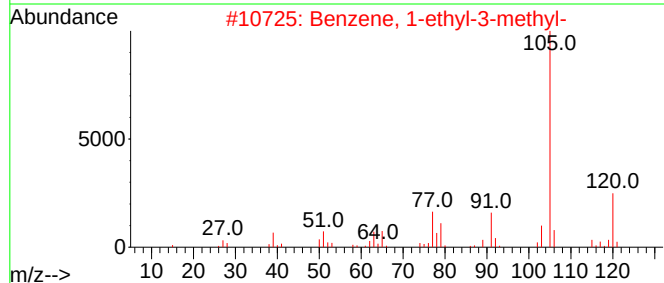
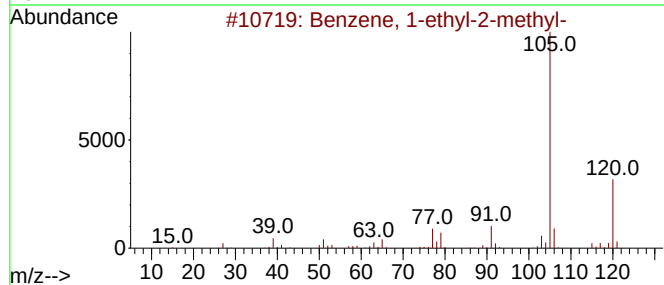
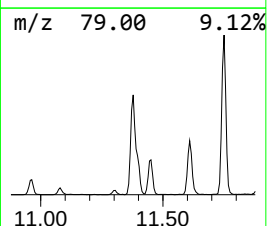
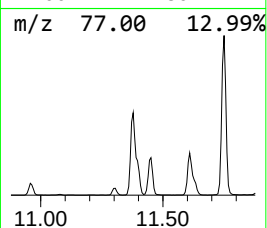
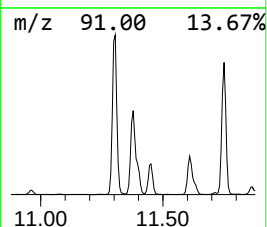
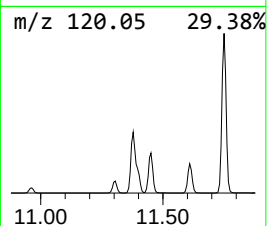
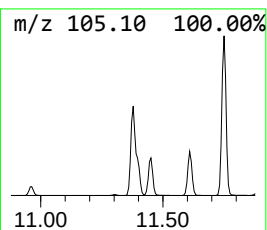
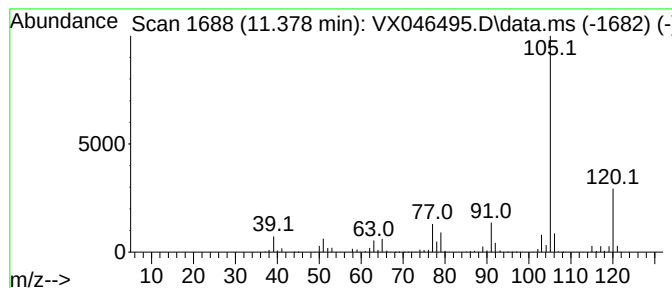
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X050525W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.378	159.45 ug/l	1750720	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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Instrument :
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ClientSampleId :
303-PPR-1

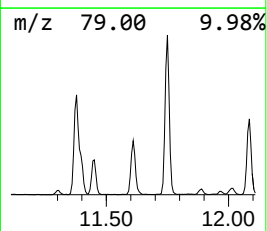
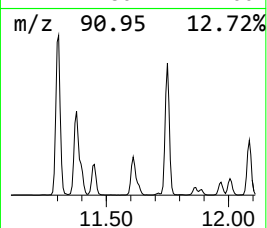
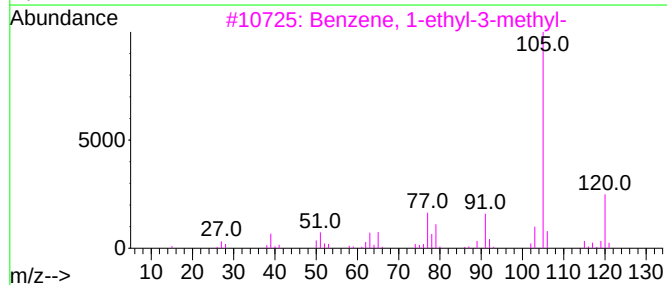
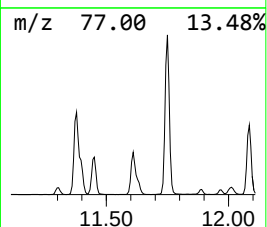
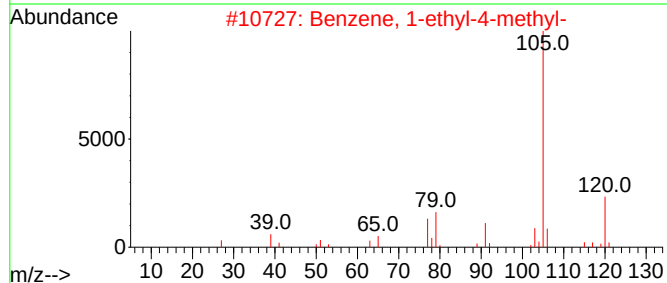
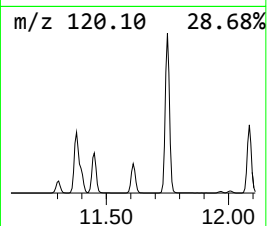
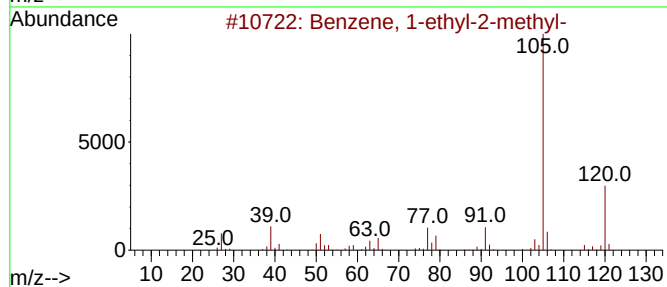
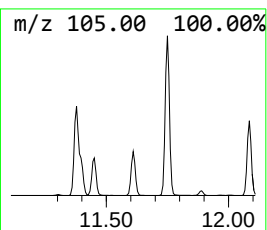
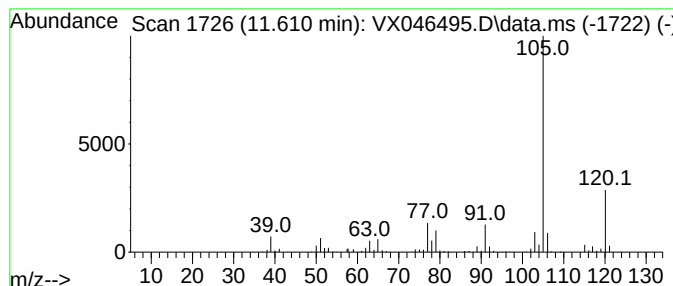
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X050525W.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1-ethyl-4-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.610	76.10 ug/l	835585	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	93
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



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ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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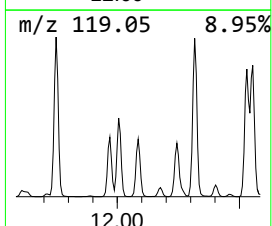
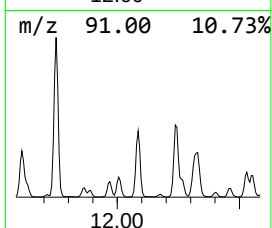
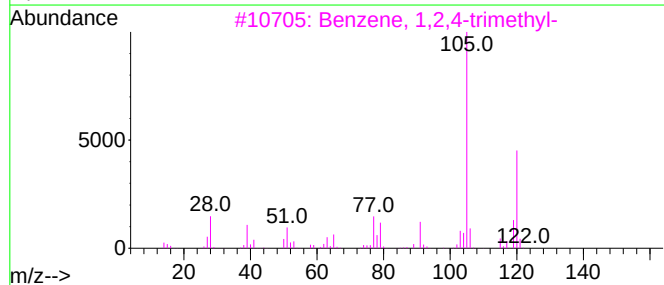
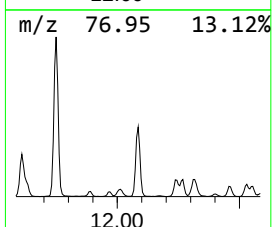
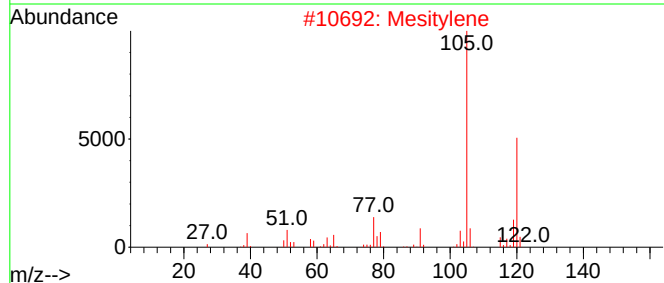
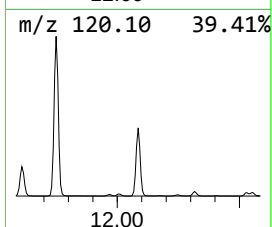
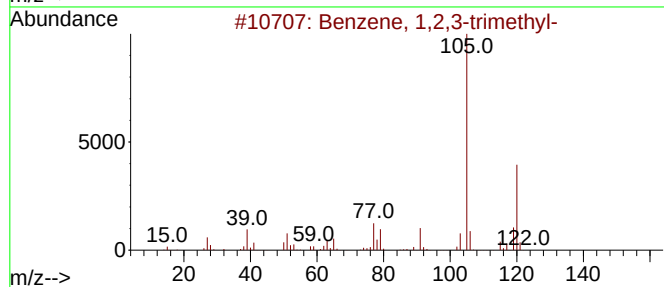
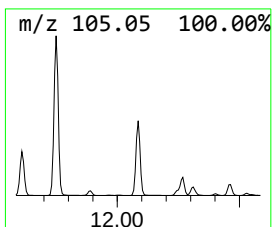
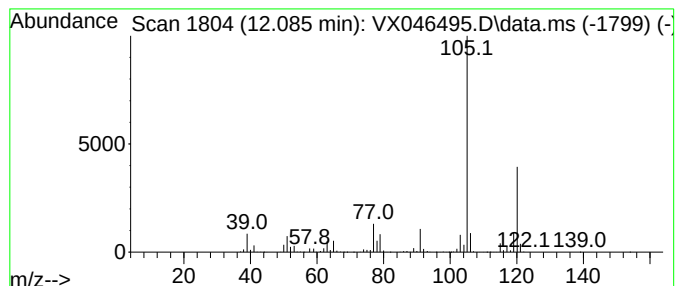
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X050525W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.085	107.01 ug/l	1174940	1,4-Dichlorobenzene-d4	12.018

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



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Instrument :
MSVOA_X
ClientSampleId :
303-PPR-1

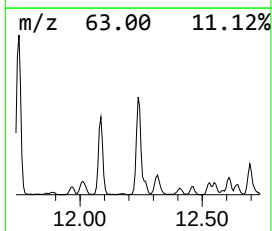
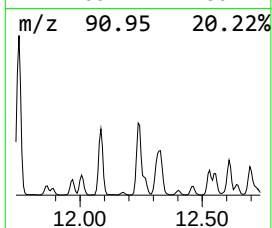
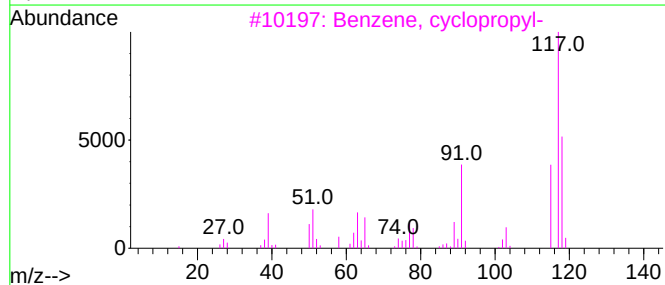
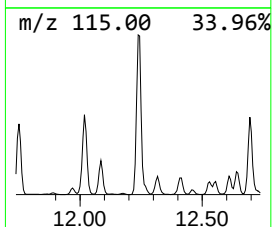
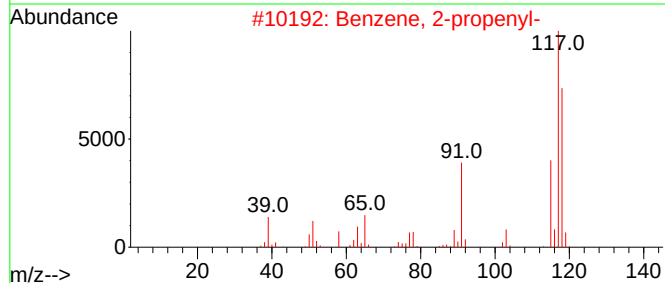
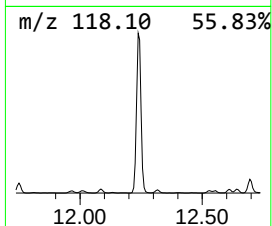
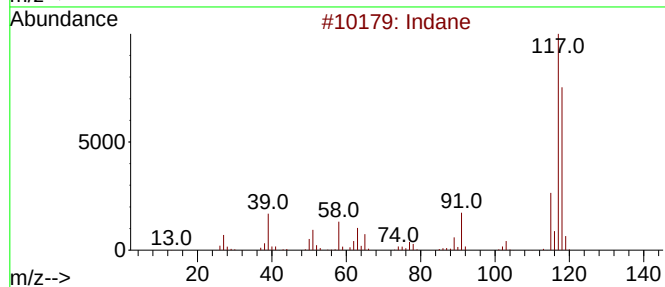
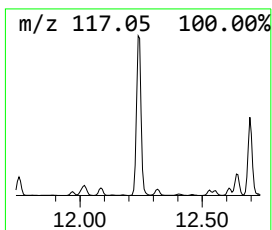
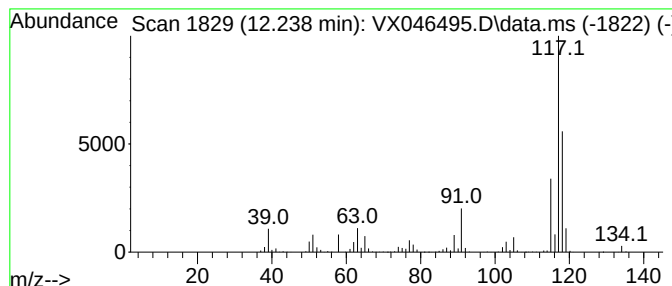
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Quant Title : SW846 8260

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

Peak Number 4 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.238	98.79 ug/l	1084620	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	87
2			Benzene, 2-propenyl-	118	C9H10	000300-57-2	81
3			Benzene, cyclopropyl-	118	C9H10	000873-49-4	81
4			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	72
5			Benzene, 1-propenyl-	118	C9H10	000637-50-3	68



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060425\
Data File : VX046495.D
Acq On : 04 Jun 2025 13:05
Operator : JC/MD
Sample : Q2169-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

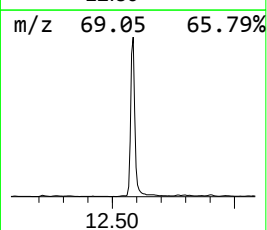
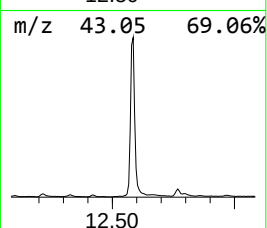
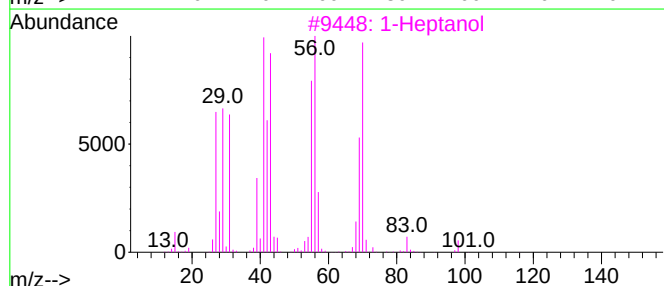
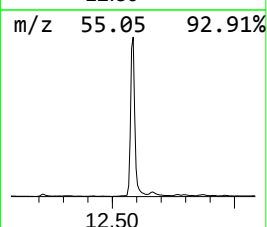
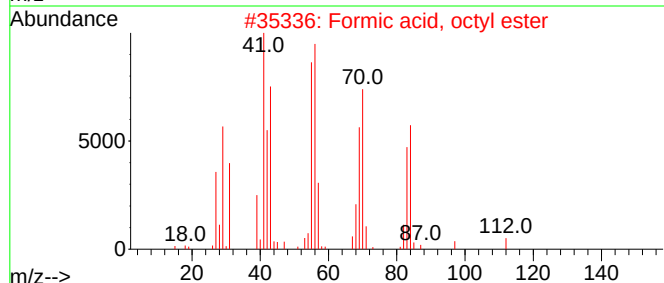
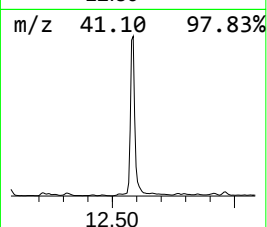
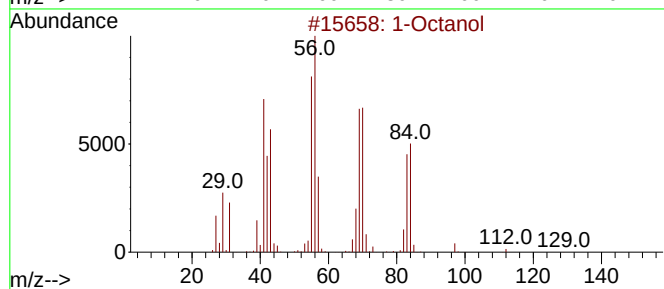
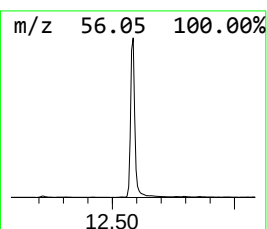
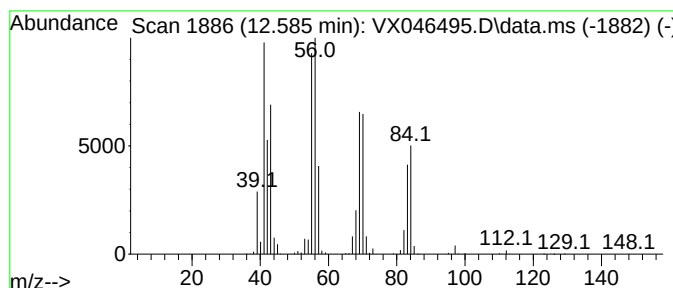
Instrument :
MSVOA_X
ClientSampleId :
303-PPR-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X050525W.M
Quant Title : SW846 8260

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

Peak Number 5 1-Octanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.585	237.16 ug/l	2603930	1,4-Dichlorobenzene-d4	12.018	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Octanol	130	C8H18O	000111-87-5	91
2	Formic acid, octyl ester	158	C9H18O2	000112-32-3	90
3	1-Heptanol	116	C7H16O	000111-70-6	50
4	1-Heptene	98	C7H14	000592-76-7	47
5	1-Hexene	84	C6H12	000592-41-6	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060425\
Data File : VX046495.D
Acq On : 04 Jun 2025 13:05
Operator : JC/MD
Sample : Q2169-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
303-PPR-1

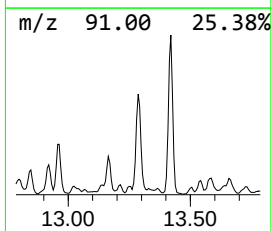
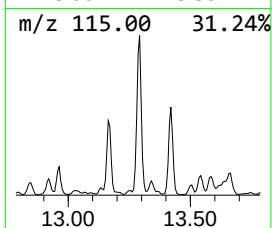
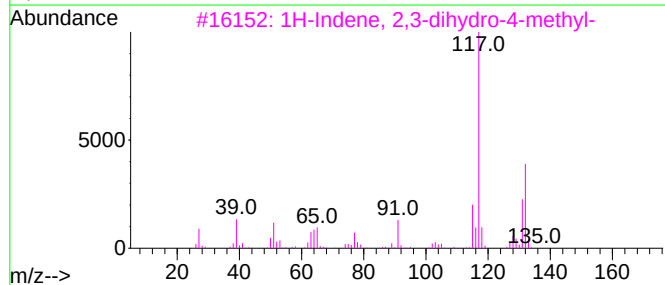
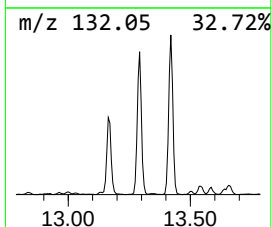
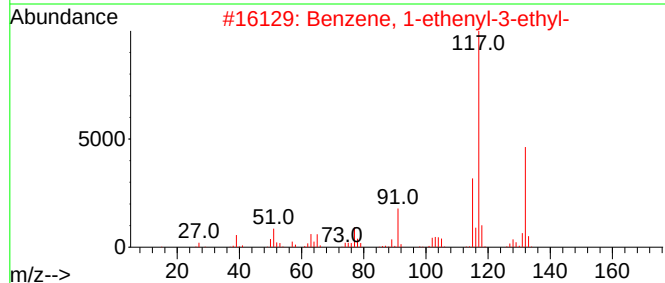
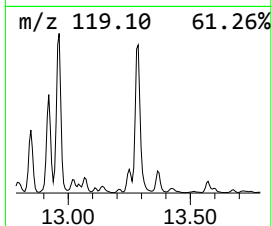
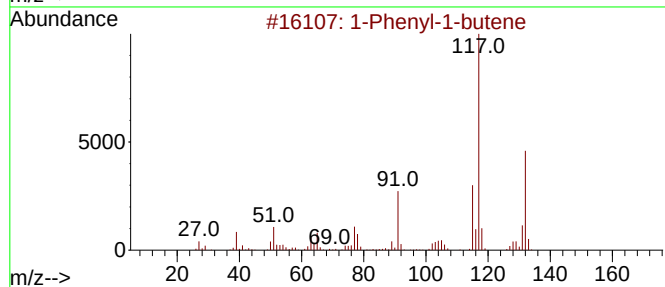
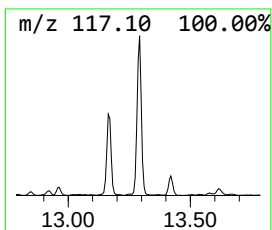
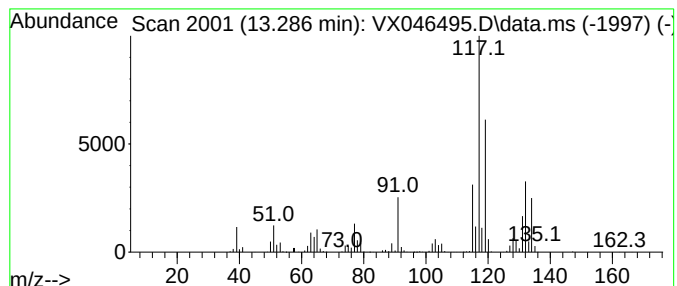
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Quant Title : SW846 8260

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

Peak Number 6 1-Phenyl-1-butene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.286	85.36 ug/l	937265	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenyl-1-butene	132	C10H12	000824-90-8	95
2			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	92
3			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	91
4			Benzene, 2-butenyl-	132	C10H12	001560-06-1	91
5			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060425\
Data File : VX046495.D
Acq On : 04 Jun 2025 13:05
Operator : JC/MD
Sample : Q2169-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
303-PPR-1

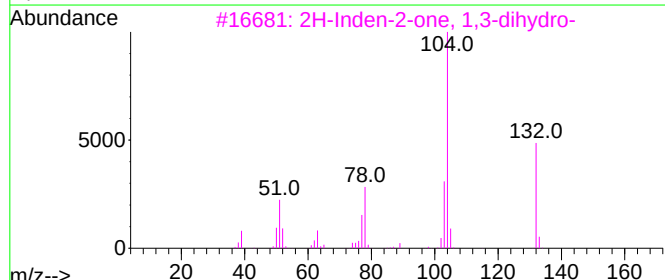
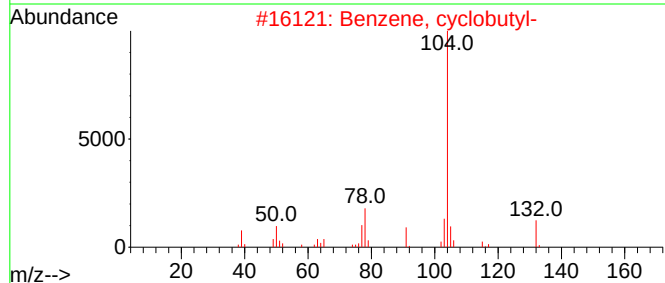
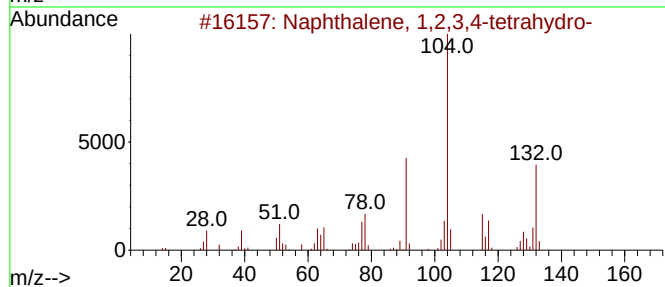
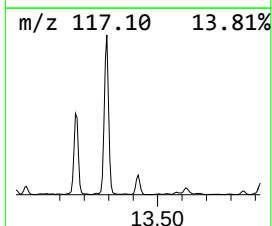
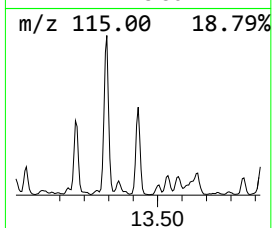
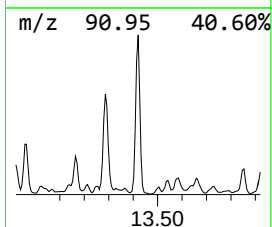
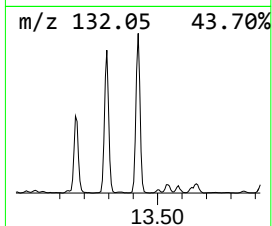
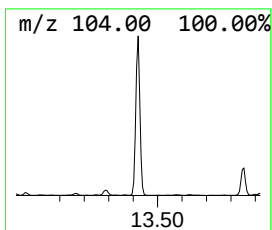
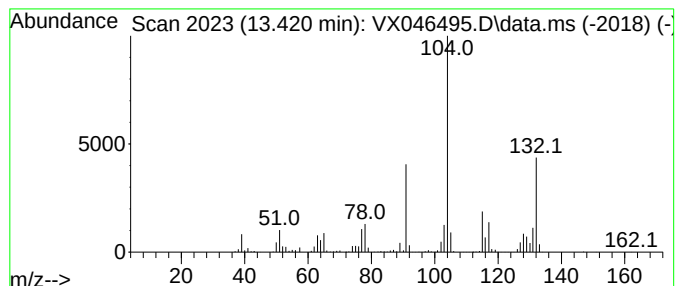
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Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.420	65.80 ug/l	722406	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	97
2			Benzene, cyclobutyl-	132	C10H12	004392-30-7	59
3			2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	58
4			Indan, epoxide	132	C9H8O	000768-22-9	52
5			2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060425\
Data File : VX046495.D
Acq On : 04 Jun 2025 13:05
Operator : JC/MD
Sample : Q2169-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
303-PPR-1

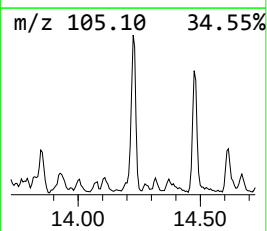
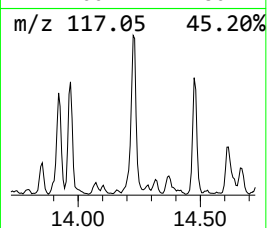
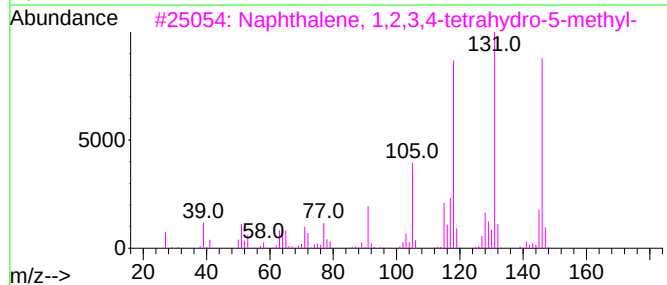
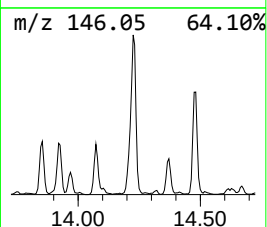
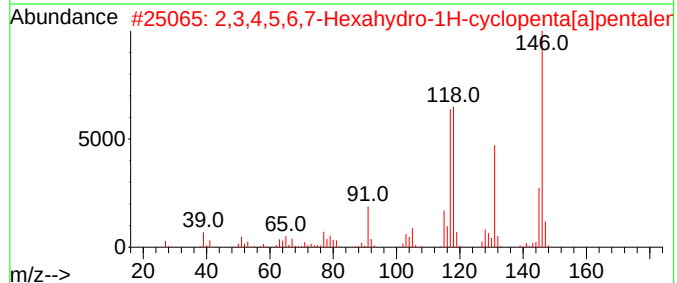
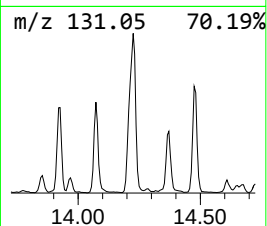
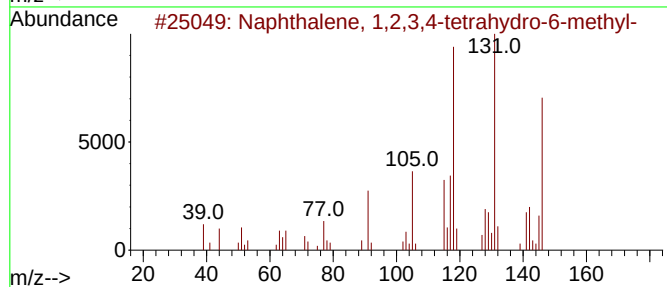
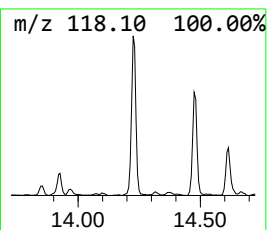
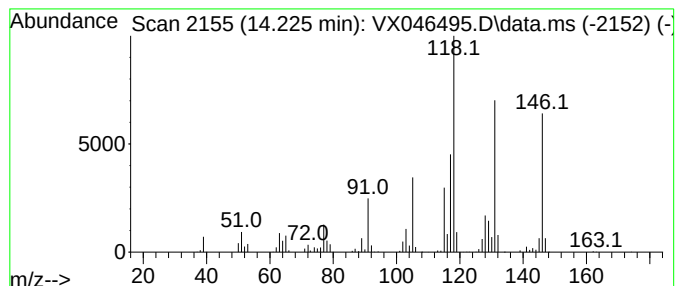
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Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.225	43.64 ug/l	479129	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	93
2			2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14	1010189-31-0	62
3			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	60
4			Benzene, (1-methylenebutyl)-	146	C11H14	005676-32-4	58
5			1H-Tetrazole, 5-phenyl-	146	C7H6N4	018039-42-4	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060425\
Data File : VX046495.D
Acq On : 04 Jun 2025 13:05
Operator : JC/MD
Sample : Q2169-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
303-PPR-1

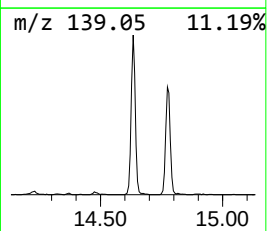
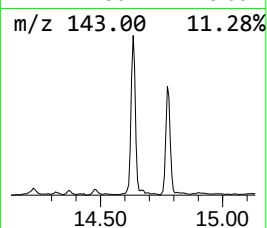
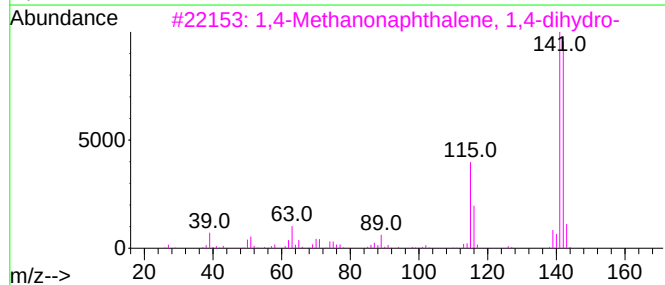
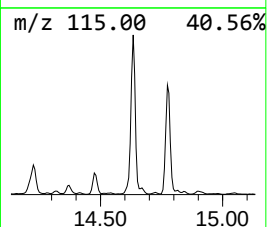
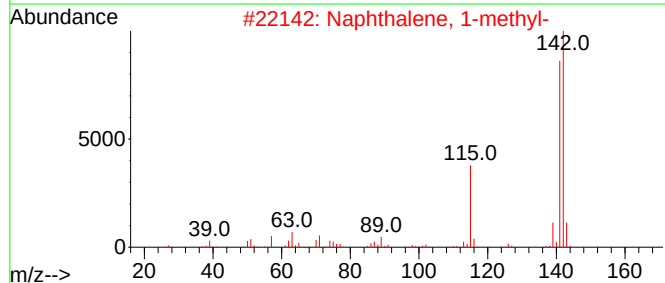
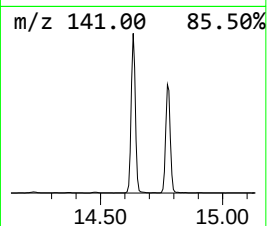
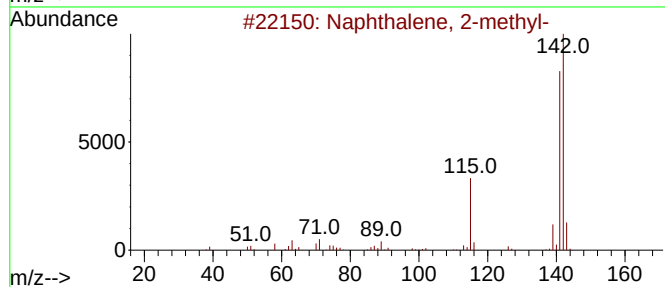
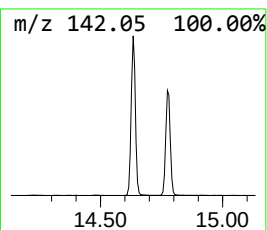
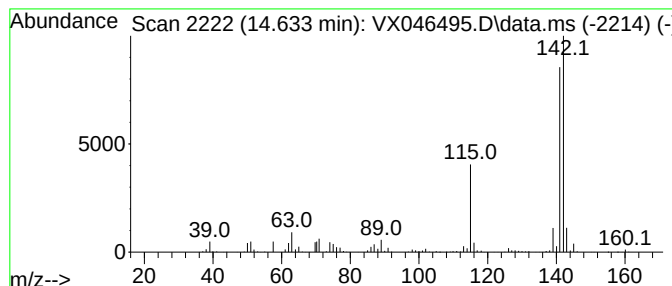
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.634	98.51 ug/l	1081590	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	97
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94
4			Benzocycloheptatriene	142	C11H10	000264-09-5	90
5			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060425\
Data File : VX046495.D
Acq On : 04 Jun 2025 13:05
Operator : JC/MD
Sample : Q2169-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
303-PPR-1

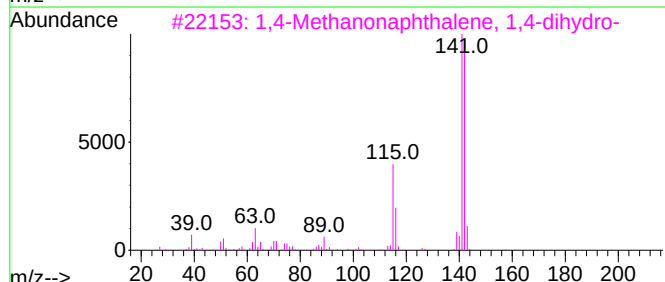
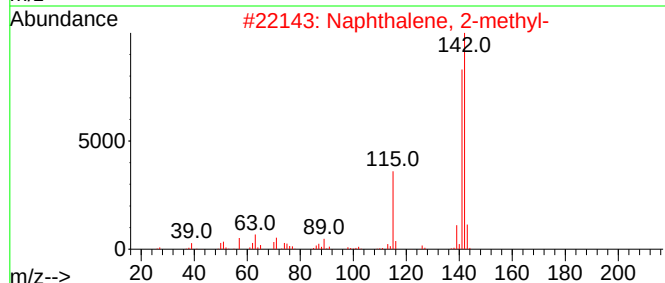
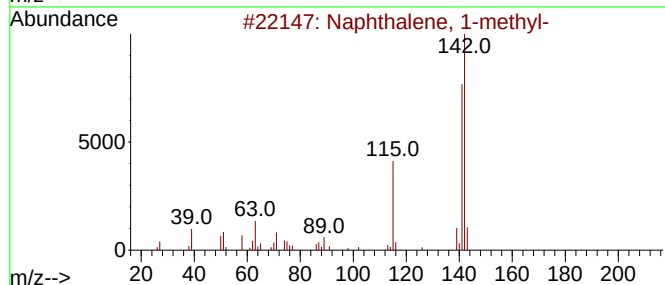
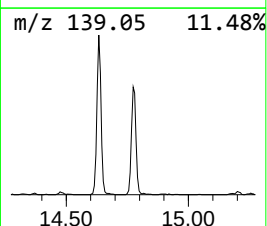
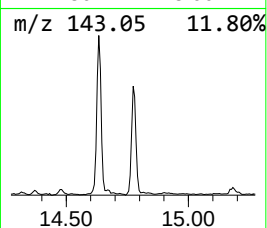
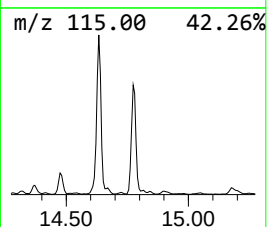
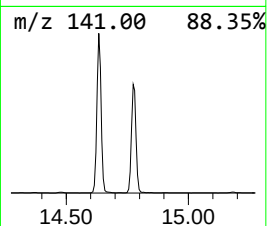
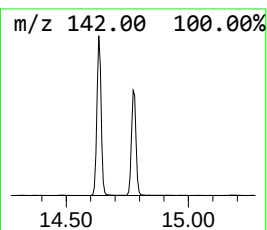
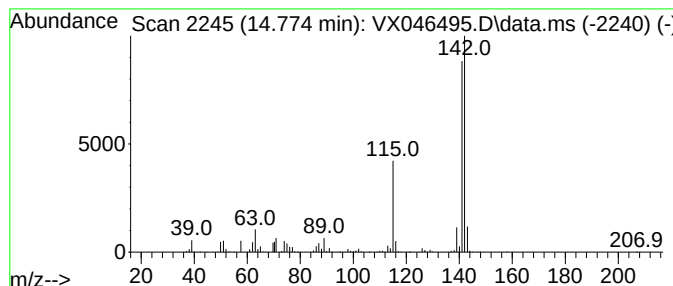
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Quant Title : SW846 8260

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.
14.774	65.69 ug/l	721209	1,4-Dichlorobenzene-d4	12.018

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	93
4			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	91
5			Benzocycloheptatriene	142	C11H10	000264-09-5	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060425\
Data File : VX046495.D
Acq On : 04 Jun 2025 13:05
Operator : JC/MD
Sample : Q2169-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
303-PPR-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X050525W.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
Benzene, 1-ethy...	11.378	159.4	ug/l	1750720	4	12.018	548979	50.0
Benzene, 1-ethy...	11.610	76.1	ug/l	835585	4	12.018	548979	50.0
Benzene, 1,2,3-...	12.085	107.0	ug/l	1174940	4	12.018	548979	50.0
Indane	12.238	98.8	ug/l	1084620	4	12.018	548979	50.0
1-Octanol	12.585	237.2	ug/l	2603930	4	12.018	548979	50.0
1-Phenyl-1-butene	13.286	85.4	ug/l	937265	4	12.018	548979	50.0
Naphthalene, 1,...	13.420	65.8	ug/l	722406	4	12.018	548979	50.0
Naphthalene, 1,...	14.225	43.6	ug/l	479129	4	12.018	548979	50.0
Naphthalene, 2-...	14.634	98.5	ug/l	1081590	4	12.018	548979	50.0
Naphthalene, 1-...	14.774	65.7	ug/l	721209	4	12.018	548979	50.0