

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX010825\
 Data File : VX044630.D
 Acq On : 08 Jan 2025 15:59
 Operator : JC/MD
 Sample : Q1036-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 NP-WS-002

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\82X010725W.M
 Title : SW846 8260

Signal : TIC: VX044630.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.246	22	26	32	rBV2	39850	44034	2.10%	0.264%
2	1.575	69	80	82	rBV3	9517	27250	1.30%	0.163%
3	2.977	299	310	320	rBV3	52031	140890	6.73%	0.845%
4	5.379	691	704	714	rBV2	120120	364086	17.40%	2.183%
5	5.544	722	731	750	rVB	200992	592586	28.32%	3.553%
6	5.946	788	797	809	rBV	143893	382666	18.29%	2.295%
7	6.239	829	845	855	rBV2	27440	93345	4.46%	0.560%
8	6.348	855	863	876	rVB4	12186	32767	1.57%	0.196%
9	6.757	921	930	944	rBV	333535	807454	38.59%	4.842%
10	7.165	990	997	1004	rBV2	10995	24871	1.19%	0.149%
11	7.354	1018	1028	1038	rBV4	16523	45709	2.18%	0.274%
12	7.976	1120	1130	1141	rVB	26476	65281	3.12%	0.391%
13	8.104	1141	1151	1154	rBV	49619	104518	4.99%	0.627%
14	8.135	1154	1156	1163	rVB	34407	52242	2.50%	0.313%
15	8.604	1225	1233	1234	rBV5	11103	20930	1.00%	0.126%
16	8.647	1234	1240	1250	rVB	664990	1057414	50.53%	6.341%
17	8.964	1287	1292	1299	rVB4	16300	36812	1.76%	0.221%
18	9.159	1319	1324	1335	rVB	29939	50180	2.40%	0.301%
19	10.049	1465	1470	1486	rBV	701215	991186	47.37%	5.944%
20	11.079	1634	1639	1650	rBV	533006	701496	33.52%	4.207%
21	11.451	1694	1700	1704	rBV	46443	62745	3.00%	0.376%
22	11.610	1721	1726	1735	rVB	130976	165426	7.91%	0.992%
23	12.018	1788	1793	1797	rBV	684545	855577	40.89%	5.130%
24	12.055	1797	1799	1801	rVV2	46052	54911	2.62%	0.329%
25	12.085	1801	1804	1809	rVB	124992	155539	7.43%	0.933%
26	12.177	1814	1819	1824	rVB	49913	62634	2.99%	0.376%
27	12.244	1824	1830	1836	rBV2	835199	1121892	53.61%	6.727%
28	12.317	1838	1842	1848	rVB	56569	77867	3.72%	0.467%
29	12.402	1851	1856	1862	rBV2	233365	309214	14.78%	1.854%
30	12.469	1862	1867	1872	rVB2	43958	65476	3.13%	0.393%
31	12.610	1885	1890	1894	rBV	817506	942238	45.03%	5.650%
32	12.640	1894	1895	1900	rVV	70827	79301	3.79%	0.476%
33	12.695	1900	1904	1912	rVV2	500183	847909	40.52%	5.085%
34	12.841	1924	1928	1933	rVB2	55427	82169	3.93%	0.493%
35	12.920	1933	1941	1945	rBV	710715	868558	41.51%	5.208%
36	12.963	1945	1948	1954	rVB2	44208	66373	3.17%	0.398%

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Peak Location: TOP

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X010725W.M
 Title : SW846 8260

37	13.030	1954	1959	1966	rBV4	49618	95962	4.59%	0.575%
38	13.091	1966	1969	1972	rBV	25983	29288	1.40%	0.176%
39	13.134	1972	1976	1980	rBV2	79239	93781	4.48%	0.562%
40	13.195	1983	1986	1991	rVB	36667	47396	2.26%	0.284%
41	13.292	1991	2002	2008	rBV2	1427863	2092632	100.00%	12.549%
42	13.347	2008	2011	2013	rBV	20572	23048	1.10%	0.138%
43	13.420	2020	2023	2030	rVB2	111906	153682	7.34%	0.922%
44	13.512	2030	2038	2040	rBV4	19543	41672	1.99%	0.250%
45	13.542	2040	2043	2045	rVV3	26042	32852	1.57%	0.197%
46	13.579	2045	2049	2054	rVV3	278822	447030	21.36%	2.681%
47	13.640	2055	2059	2060	rVV	143761	187634	8.97%	1.125%
48	13.664	2060	2063	2073	rVV2	239238	415150	19.84%	2.489%
49	13.750	2073	2077	2080	rVV2	22259	25480	1.22%	0.153%
50	13.792	2080	2084	2089	rVB2	63376	86955	4.16%	0.521%
51	13.853	2089	2094	2099	rVB4	38503	56679	2.71%	0.340%
52	13.926	2099	2106	2110	rBV2	112520	170197	8.13%	1.021%
53	13.969	2110	2113	2117	rVB2	96048	109230	5.22%	0.655%
54	14.073	2125	2130	2133	rBV2	53111	65263	3.12%	0.391%
55	14.213	2148	2153	2160	rBV	207153	276208	13.20%	1.656%
56	14.317	2167	2170	2175	rVB	106291	124896	5.97%	0.749%
57	14.371	2175	2179	2183	rVV	114146	139566	6.67%	0.837%
58	14.420	2183	2187	2192	rVB	24591	35761	1.71%	0.214%
59	14.481	2192	2197	2201	rBV2	106389	152159	7.27%	0.912%
60	14.615	2215	2219	2221	rBV3	28972	39203	1.87%	0.235%
61	14.670	2225	2228	2232	rVB	29464	38783	1.85%	0.233%
62	14.786	2243	2247	2249	rVV3	16498	27086	1.29%	0.162%
63	14.816	2249	2252	2255	rVV2	29224	44456	2.12%	0.267%
64	15.048	2286	2290	2298	rBV2	14136	21044	1.01%	0.126%
65	15.182	2306	2312	2319	rBV3	50151	100691	4.81%	0.604%
66	15.524	2364	2368	2373	rVB4	13773	24204	1.16%	0.145%
67	15.804	2410	2414	2419	rBV3	14684	24766	1.18%	0.149%

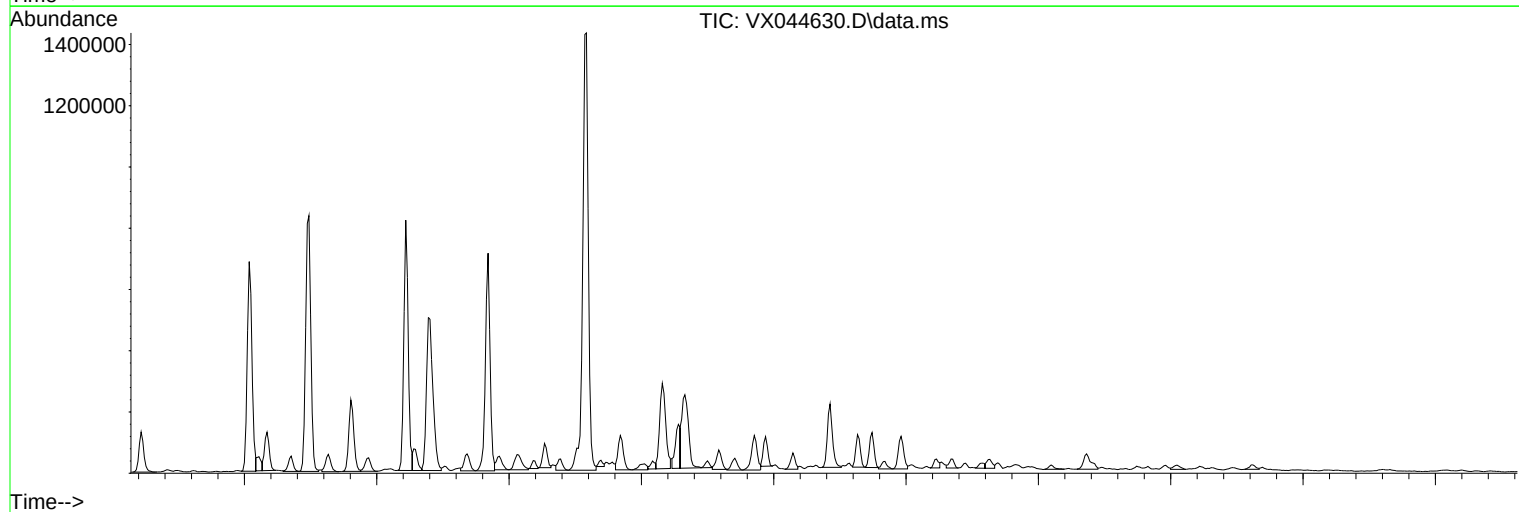
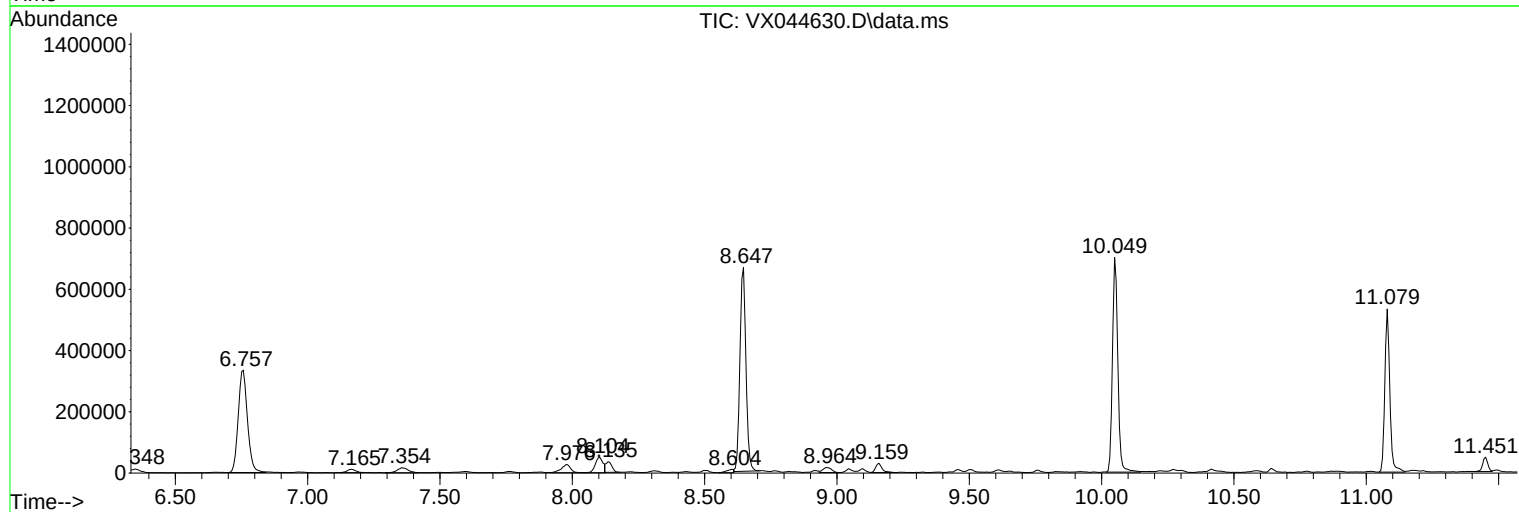
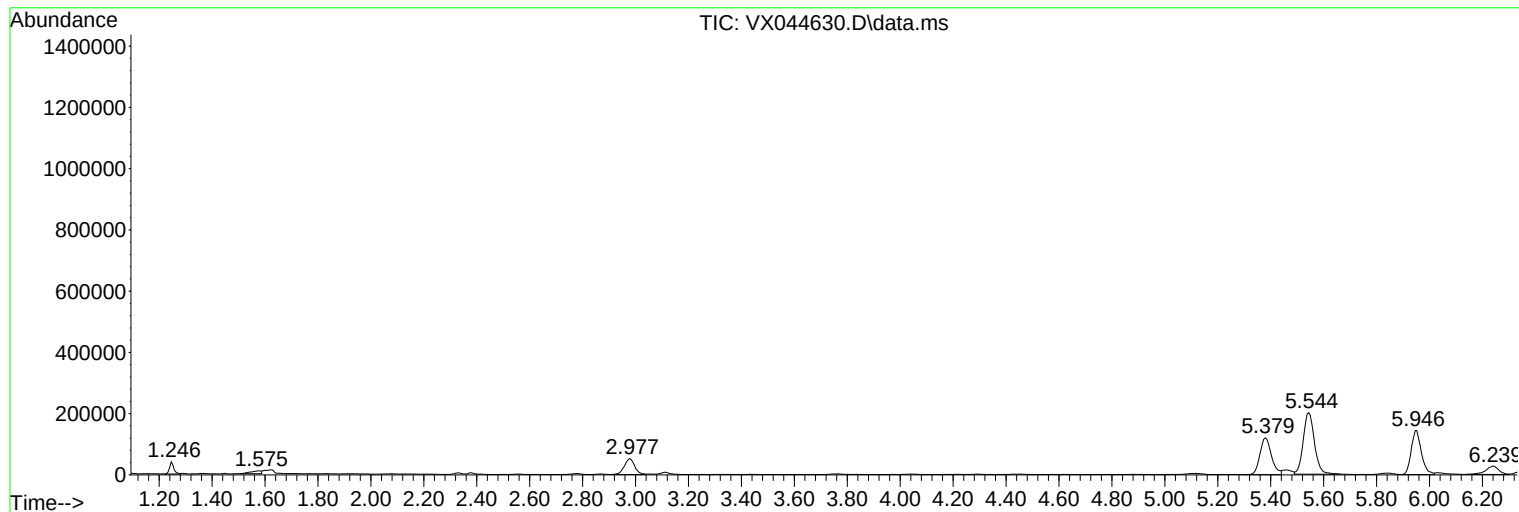
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ClientSampleId :
NP-WS-002

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X010725W.M
Quant Title : SW846 8260

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



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NP-WS-002

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X010725W.M
Quant Title : SW846 8260

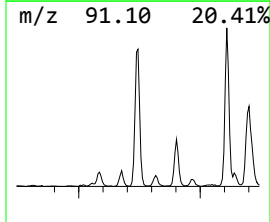
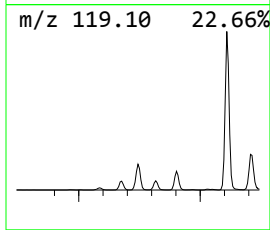
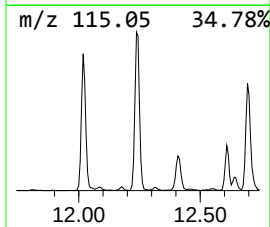
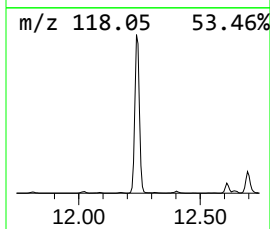
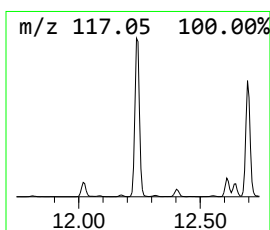
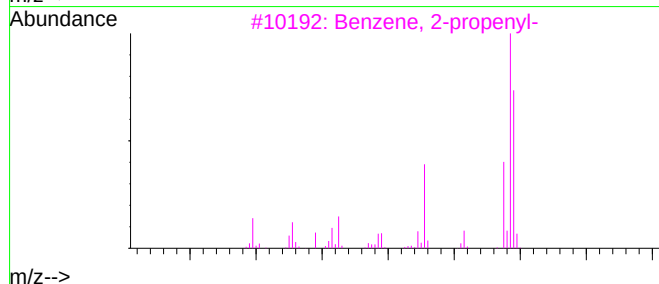
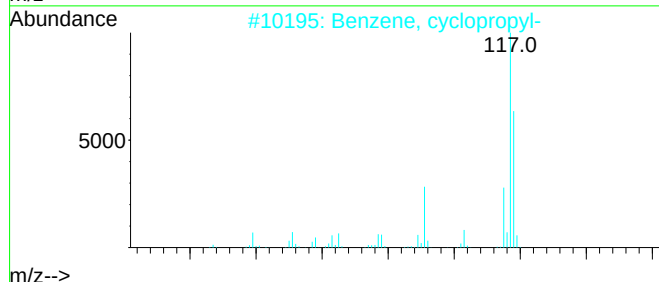
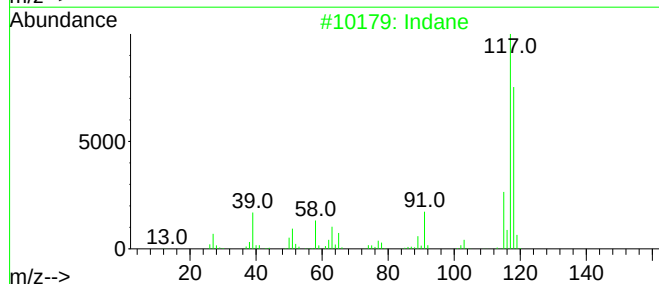
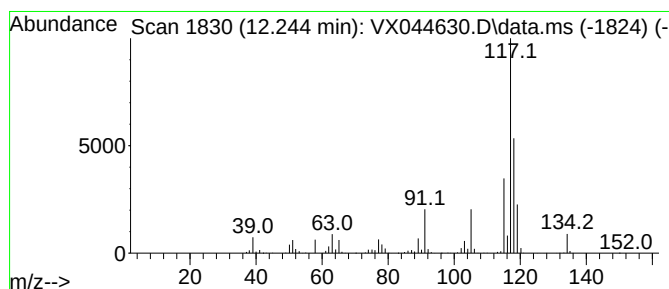
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 1 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	65.56 ug/l	1121890	1,4-Dichlorobenzene-d4	12.018

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	70
2	Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	70
4	Benzene, 1-propenyl-	118	C9H10	000637-50-3	62
5	Benzeneethanol, .beta.-ethenyl-	148	C10H12O	006052-63-7	59



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MSVOA_X
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NP-WS-002

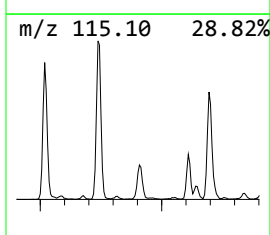
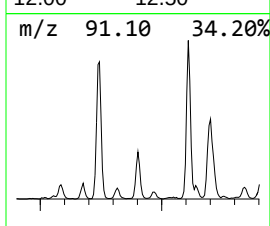
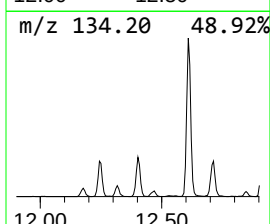
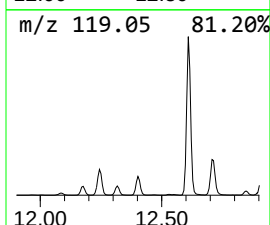
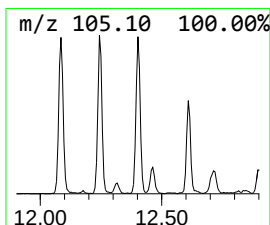
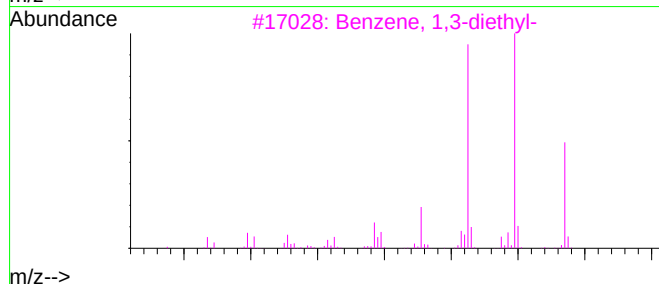
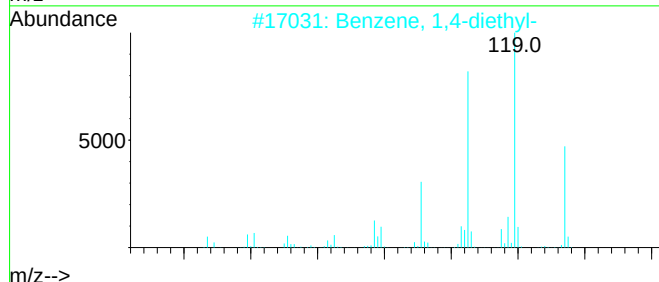
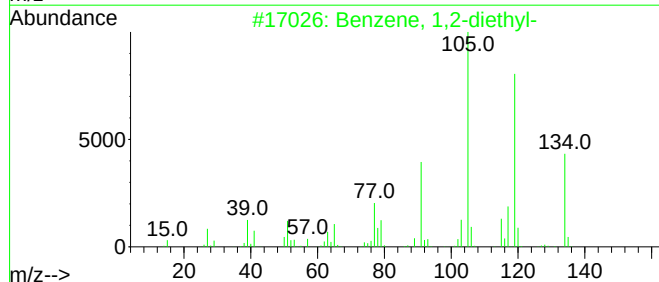
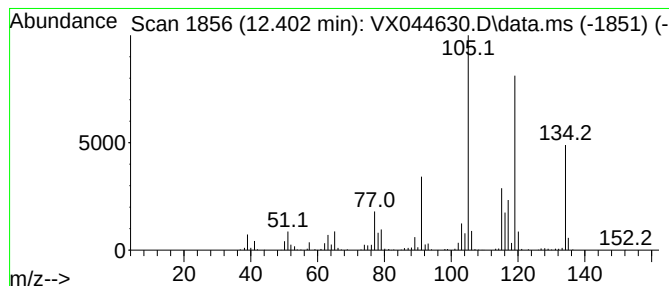
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X010725W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 1,2-diethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.402	18.07 ug/l	309214	1,4-Dichlorobenzene-d4	12.018	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	94
2	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	81
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	74
4	2,6-Dimethyl-1,3,5,7-octatetraen...	134	C10H14	000460-01-5	74
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	58



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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X010725W.M
Quant Title : SW846 8260

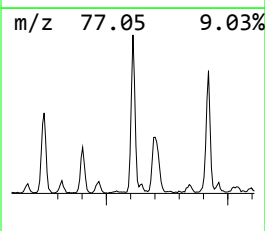
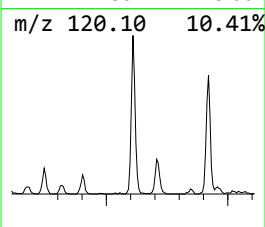
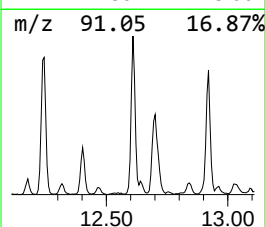
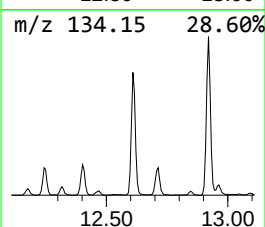
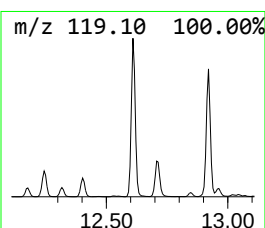
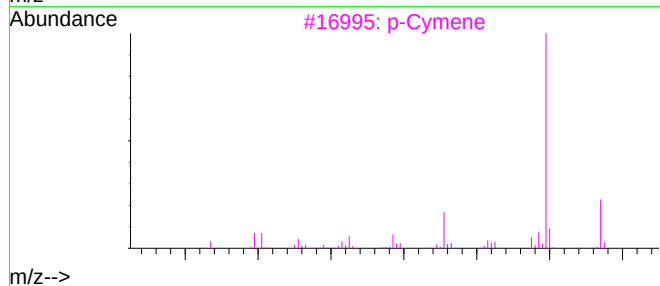
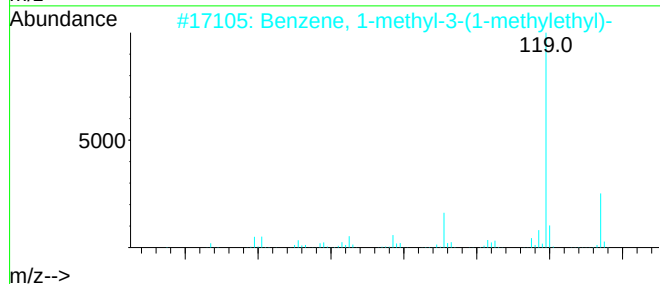
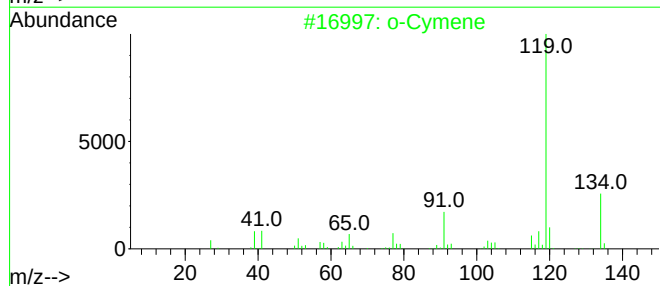
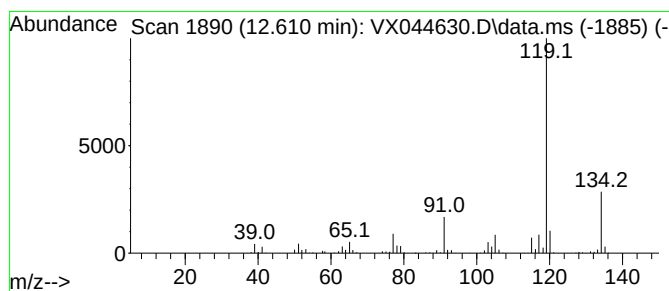
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 3 o-Cymene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.610	55.06 ug/l	942238	1,4-Dichlorobenzene-d4	12.018

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	97
2	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
3	p-Cymene	134	C10H14	000099-87-6	95
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



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Instrument :
MSVOA_X
ClientSampleId :
NP-WS-002

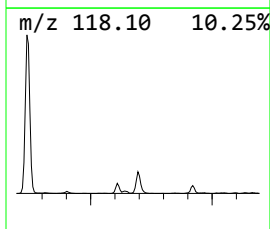
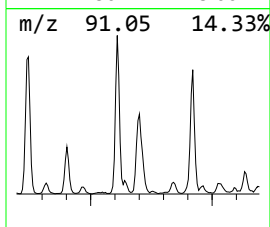
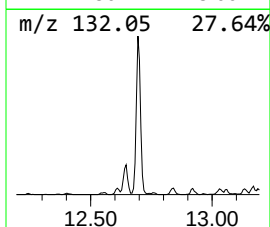
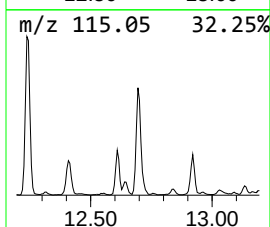
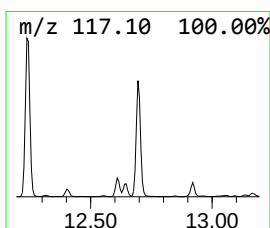
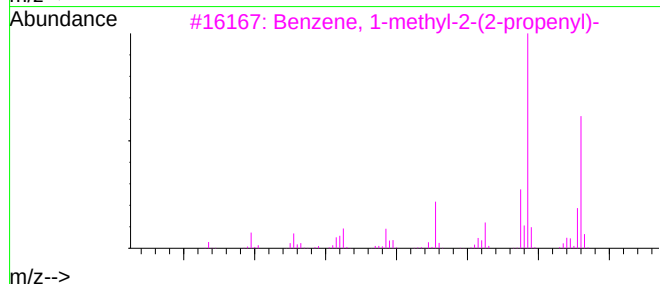
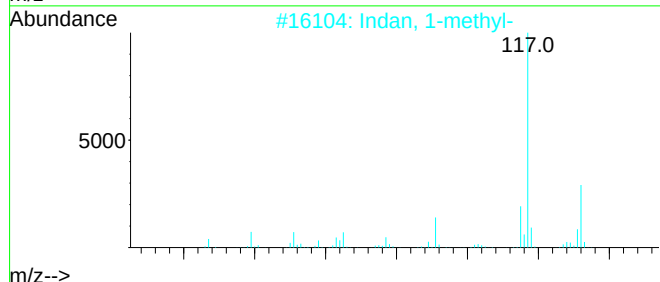
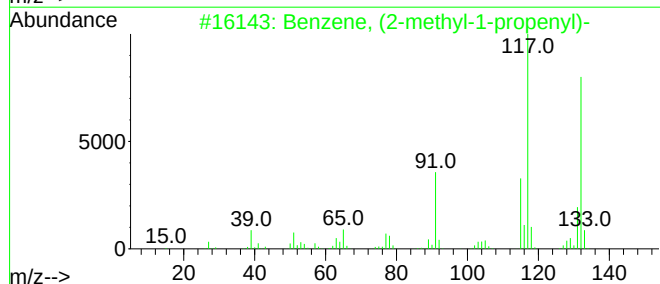
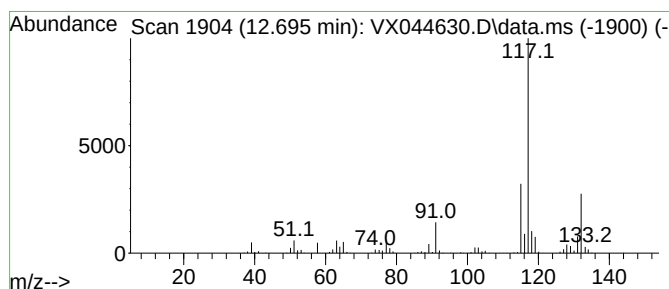
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X010725W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, (2-methyl-1-propen... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.695	49.55 ug/l	847909	1,4-Dichlorobenzene-d4	12.018	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90
2	Indan, 1-methyl-	132	C10H12	000767-58-8	87
3	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	83
4	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	83
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX010825\
Data File : VX044630.D
Acq On : 08 Jan 2025 15:59
Operator : JC/MD
Sample : Q1036-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
NP-WS-002

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X010725W.M
Quant Title : SW846 8260

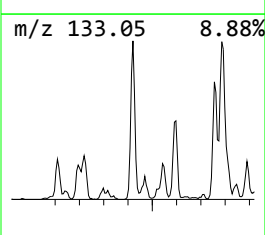
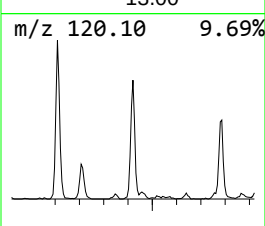
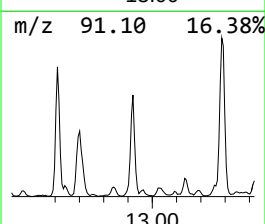
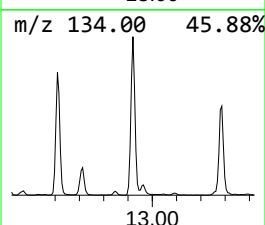
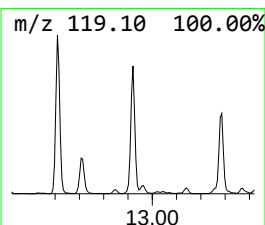
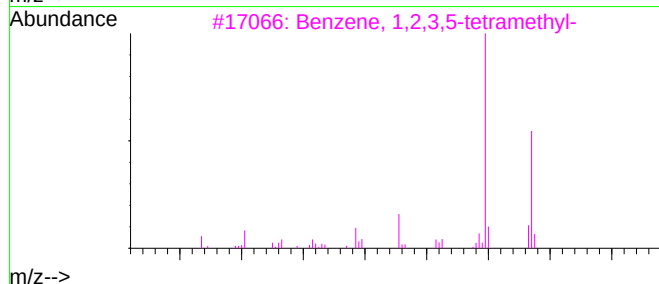
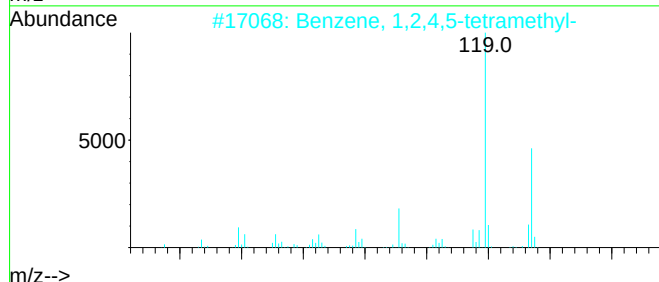
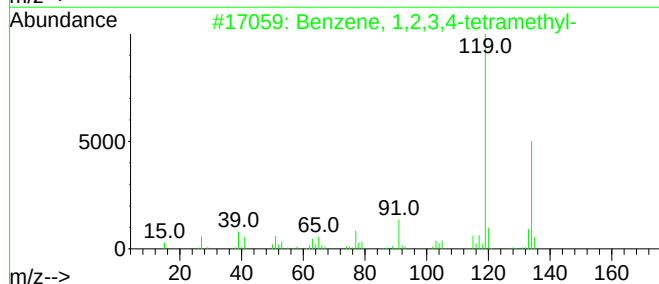
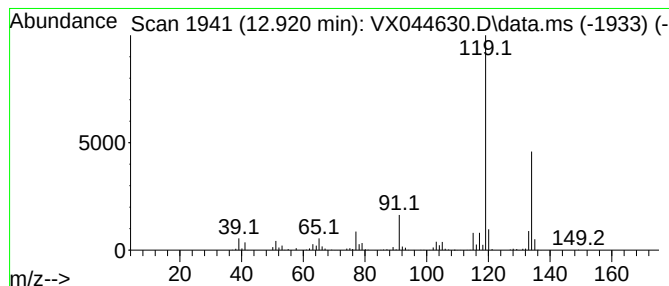
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.920	50.76 ug/l	868558	1,4-Dichlorobenzene-d4	12.018

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-2,4-dimethyl-	134	C10H14	000874-41-9	95
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX010825\
Data File : VX044630.D
Acq On : 08 Jan 2025 15:59
Operator : JC/MD
Sample : Q1036-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
NP-WS-002

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X010725W.M
Quant Title : SW846 8260

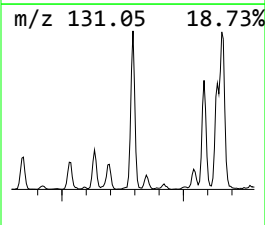
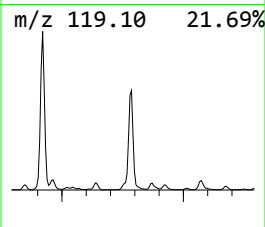
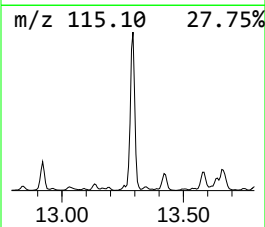
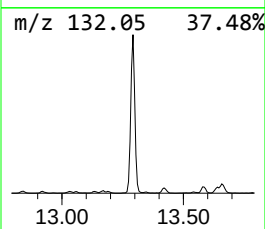
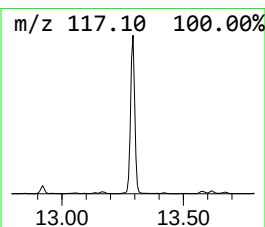
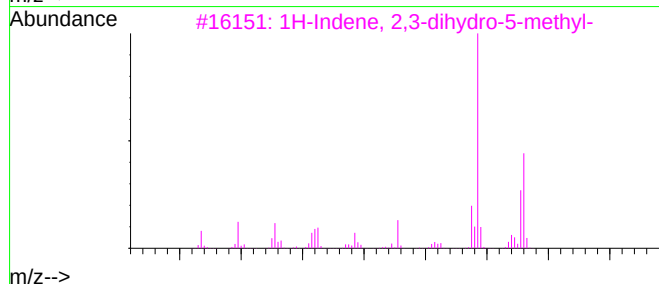
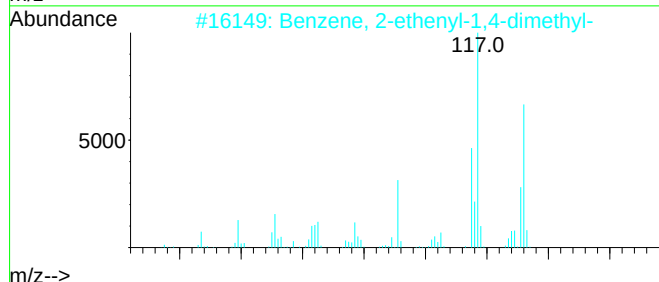
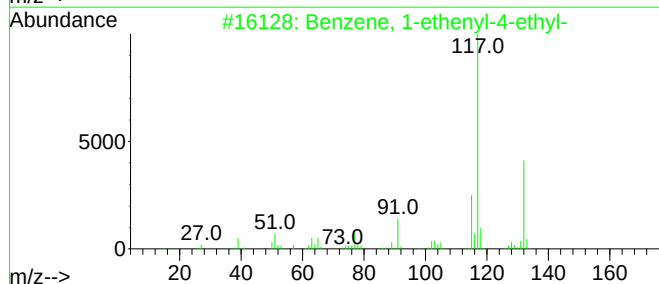
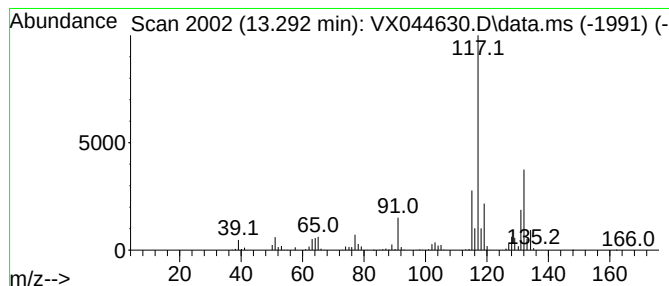
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	122.29 ug/l	2092630	1,4-Dichlorobenzene-d4	12.018

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	90
3	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	89
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
5	Indan, 1-methyl-	132	C10H12	000767-58-8	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX010825\
Data File : VX044630.D
Acq On : 08 Jan 2025 15:59
Operator : JC/MD
Sample : Q1036-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
NP-WS-002

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X010725W.M
Quant Title : SW846 8260

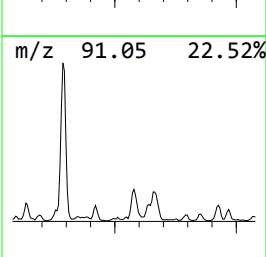
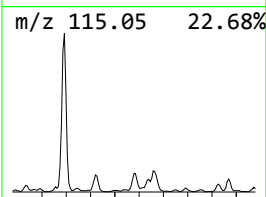
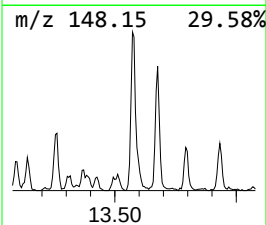
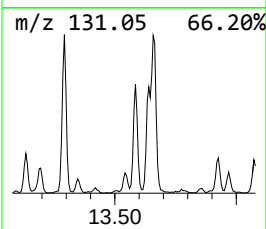
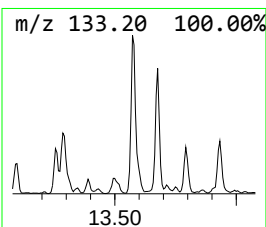
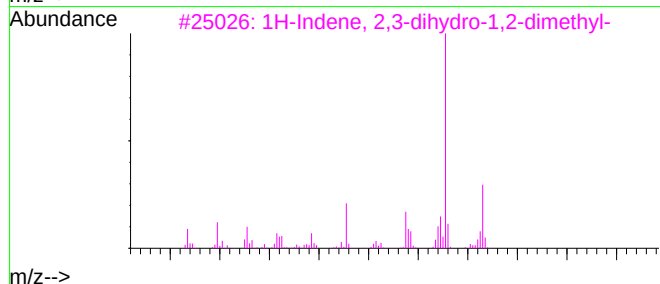
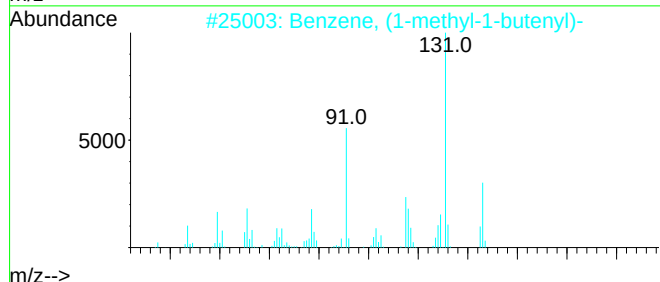
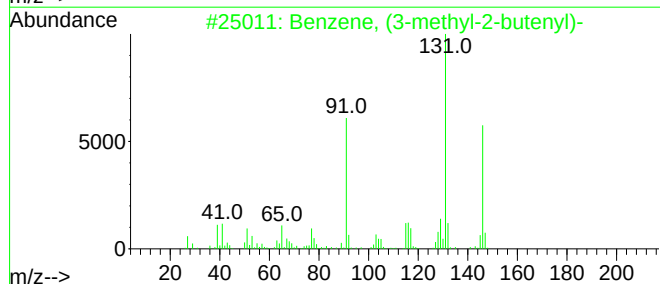
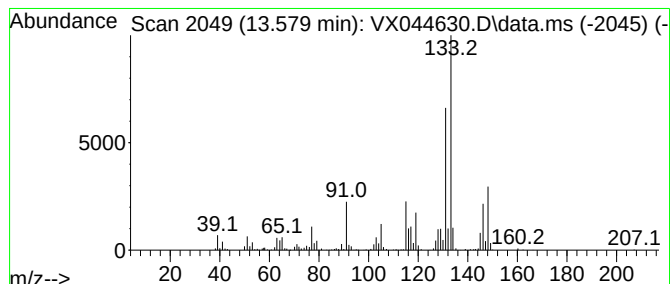
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, (3-methyl-2-butenyl)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.579	26.12 ug/l	447030	1,4-Dichlorobenzene-d4	12.018

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	86
2	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	84
3	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	62
4	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	53
5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX010825\
Data File : VX044630.D
Acq On : 08 Jan 2025 15:59
Operator : JC/MD
Sample : Q1036-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
NP-WS-002

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X010725W.M
Quant Title : SW846 8260

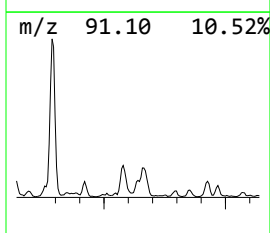
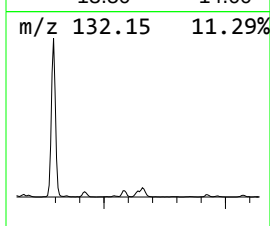
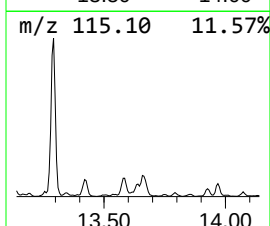
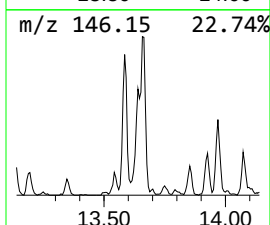
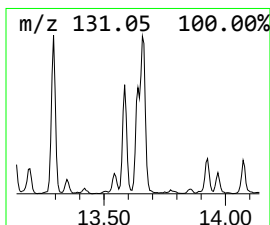
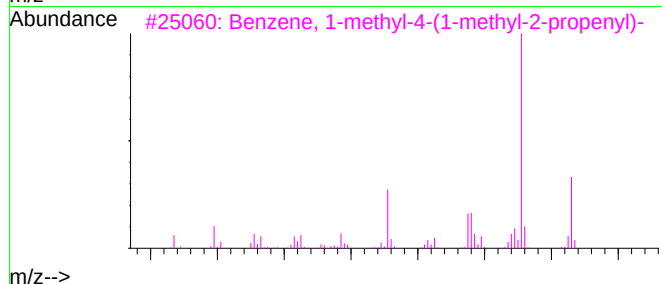
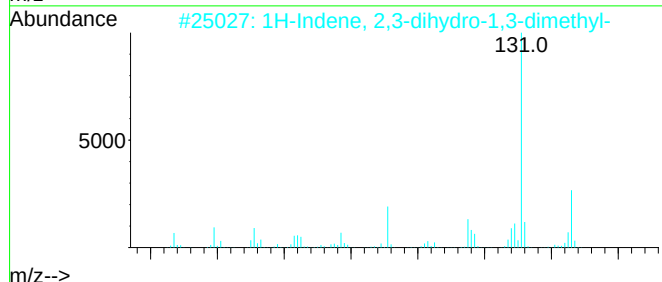
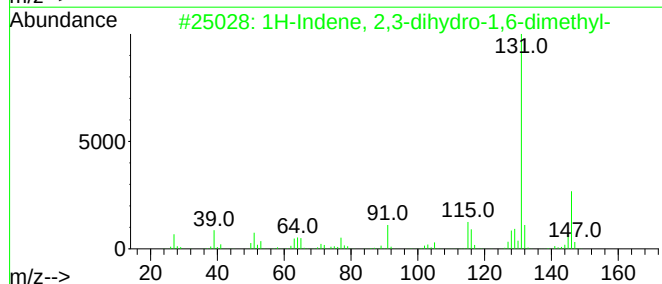
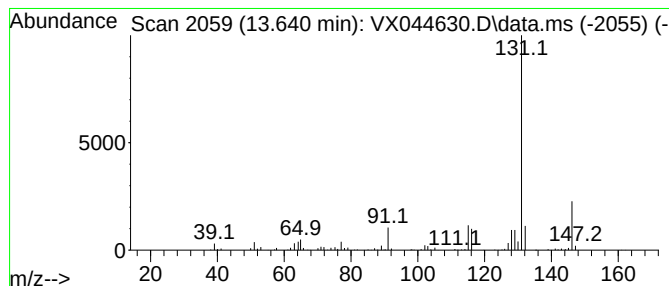
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 8 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.640	10.97 ug/l	187634	1,4-Dichlorobenzene-d4	12.018

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	91
2	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
3	Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	90
4	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	90
5	Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX010825\
Data File : VX044630.D
Acq On : 08 Jan 2025 15:59
Operator : JC/MD
Sample : Q1036-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
NP-WS-002

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X010725W.M
Quant Title : SW846 8260

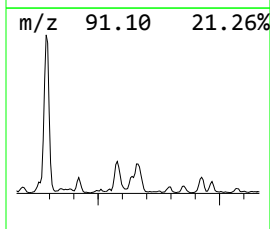
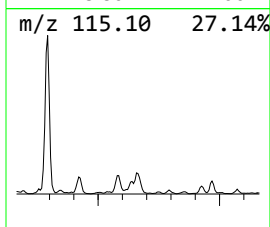
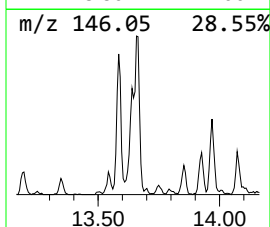
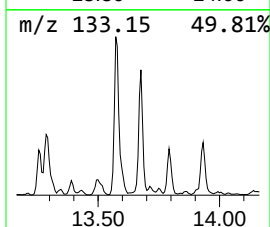
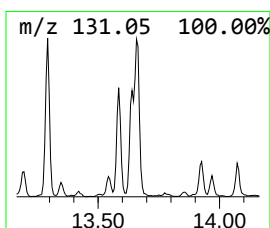
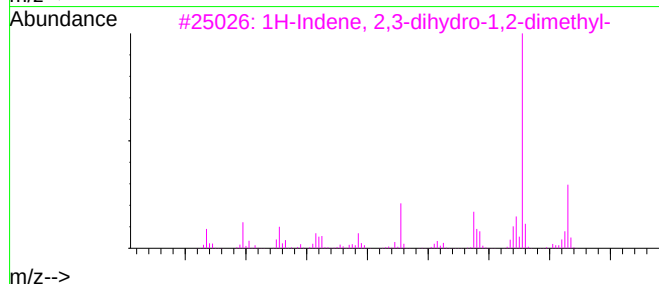
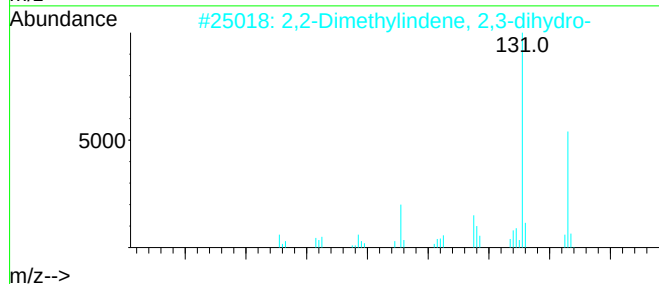
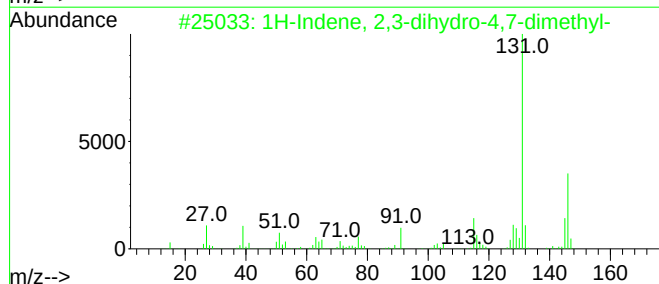
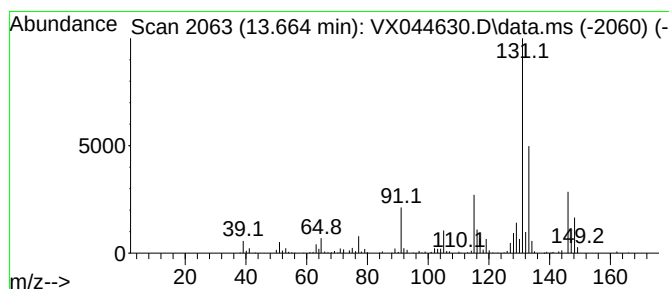
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 9 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.664	24.26 ug/l	415150	1,4-Dichlorobenzene-d4	12.018

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	70
2	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	70
3	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	70
4	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	70
5	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX010825\
Data File : VX044630.D
Acq On : 08 Jan 2025 15:59
Operator : JC/MD
Sample : Q1036-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
NP-WS-002

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X010725W.M
Quant Title : SW846 8260

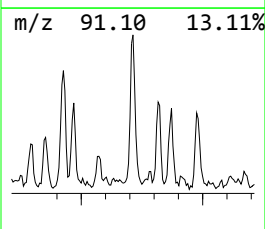
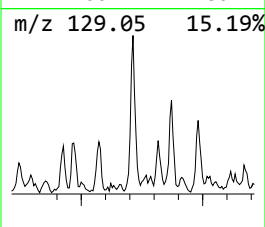
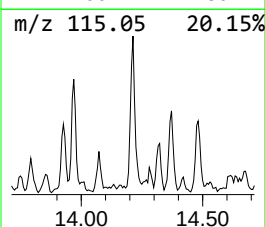
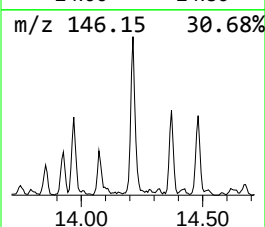
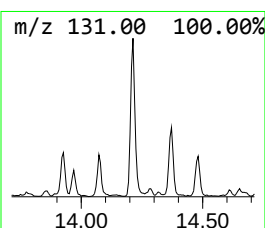
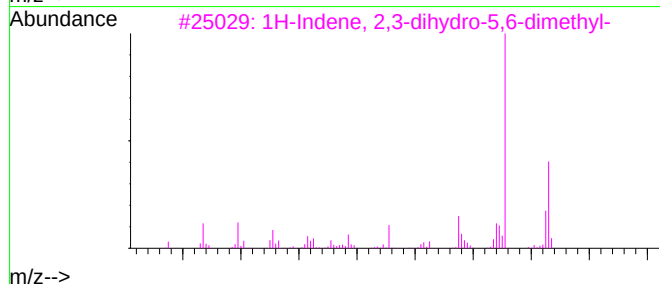
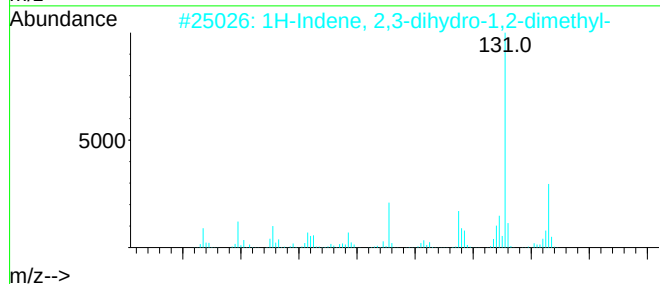
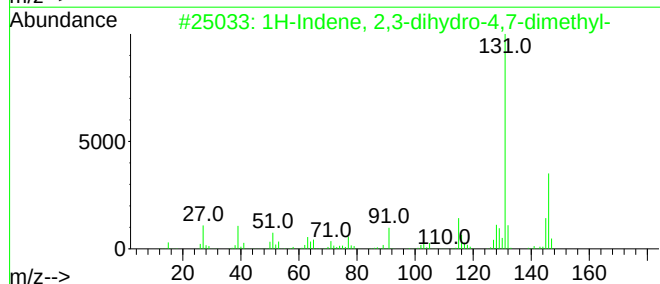
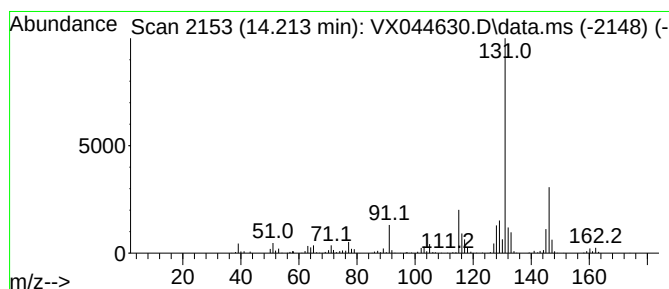
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 10 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.213	16.14 ug/l	276208	1,4-Dichlorobenzene-d4	12.018

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
2	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	94
3	1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	91
4	Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	90
5	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX010825\
Data File : VX044630.D
Acq On : 08 Jan 2025 15:59
Operator : JC/MD
Sample : Q1036-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
NP-WS-002

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X010725W.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
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Benzene, 1,2-di...	12.402	18.1	ug/l	309214	4	12.018	855577	50.0
o-Cymene	12.610	55.1	ug/l	942238	4	12.018	855577	50.0
Benzene, (2-met...	12.695	49.5	ug/l	847909	4	12.018	855577	50.0
Benzene, 1,2,3,...	12.920	50.8	ug/l	868558	4	12.018	855577	50.0
Benzene, 1-ethe...	13.292	122.3	ug/l	2092630	4	12.018	855577	50.0
Benzene, (3-met...	13.579	26.1	ug/l	447030	4	12.018	855577	50.0
1H-Indene, 2,3-...	13.640	11.0	ug/l	187634	4	12.018	855577	50.0
1H-Indene, 2,3-...	13.664	24.3	ug/l	415150	4	12.018	855577	50.0
1H-Indene, 2,3-...	14.213	16.1	ug/l	276208	4	12.018	855577	50.0