

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD031822\
 Data File : VD072324.D
 Acq On : 19 Mar 2022 08:23
 Operator : VA/SY
 Sample : N1972-02
 Misc : 5.06G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 G-4-7-17-18

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_D\Method\82D031122S.M
 Title : SW846 8260

Signal : TIC: VD072324.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.914	1011	1019	1024	rBV3	15329	39213	7.16%	1.595%
2	7.973	1024	1029	1041	rVB2	23490	55060	10.05%	2.239%
3	8.350	1083	1093	1101	rVB4	45376	127322	23.24%	5.179%
4	8.861	1170	1180	1187	rVB2	36329	70646	12.89%	2.873%
5	10.338	1422	1431	1437	rBV	54674	99464	18.15%	4.046%
6	10.514	1454	1461	1466	rVB4	5109	8249	1.51%	0.336%
7	11.238	1580	1584	1588	rBV3	3535	7118	1.30%	0.290%
8	11.638	1646	1652	1663	rBV2	59395	107853	19.68%	4.387%
9	11.738	1663	1669	1676	rBV2	51126	88465	16.15%	3.598%
10	11.849	1681	1688	1696	rVV3	24690	51373	9.38%	2.090%
11	12.173	1737	1743	1748	rBV6	5788	9896	1.81%	0.403%
12	12.467	1788	1793	1799	rBV2	19519	31695	5.78%	1.289%
13	12.626	1813	1820	1827	rVB2	53514	85970	15.69%	3.497%
14	12.726	1833	1837	1843	rVB7	7982	13003	2.37%	0.529%
15	12.808	1843	1851	1856	rBV4	5308	11427	2.09%	0.465%
16	12.914	1862	1869	1871	rBV3	11954	21568	3.94%	0.877%
17	12.949	1871	1875	1881	rVB6	13387	24944	4.55%	1.015%
18	13.114	1897	1903	1908	rVB	17924	31246	5.70%	1.271%
19	13.261	1921	1928	1934	rBV2	43832	72130	13.16%	2.934%
20	13.508	1965	1970	1973	rBV3	7329	11155	2.04%	0.454%
21	13.567	1973	1980	1982	rBV	70989	122140	22.29%	4.968%
22	13.767	2001	2014	2025	rBV	315279	547897	100.00%	22.285%
23	13.885	2030	2034	2039	rVB5	5788	9039	1.65%	0.368%
24	13.949	2039	2045	2051	rBV5	7689	16168	2.95%	0.658%
25	14.043	2059	2061	2065	rVB3	6406	7958	1.45%	0.324%
26	14.108	2065	2072	2076	rBV3	13436	21792	3.98%	0.886%
27	14.214	2087	2090	2097	rVB4	15034	22339	4.08%	0.909%
28	14.343	2108	2112	2116	rVB4	4872	8325	1.52%	0.339%
29	14.420	2121	2125	2130	rVV4	8509	18885	3.45%	0.768%
30	14.461	2130	2132	2137	rVV4	13096	19963	3.64%	0.812%
31	14.537	2140	2145	2150	rVV6	7416	16871	3.08%	0.686%
32	14.696	2168	2172	2177	rVB4	13040	19360	3.53%	0.787%
33	14.808	2184	2191	2198	rBV6	24383	54225	9.90%	2.206%
34	14.879	2198	2203	2210	rVV4	12397	21079	3.85%	0.857%
35	14.967	2213	2218	2224	rVB5	19192	37076	6.77%	1.508%
36	15.073	2233	2236	2240	rVB6	4916	5532	1.01%	0.225%

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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_D\Method\82D031122S.M
Title : SW846 8260

37	15.202	2255	2258	2265	rVB5	5072	6331	1.16%	0.258%
38	15.384	2281	2289	2297	rBV	261672	476068	86.89%	19.363%
39	16.473	2469	2474	2480	rVB3	13794	29947	5.47%	1.218%
40	16.684	2502	2510	2515	rBV5	14121	29828	5.44%	1.213%

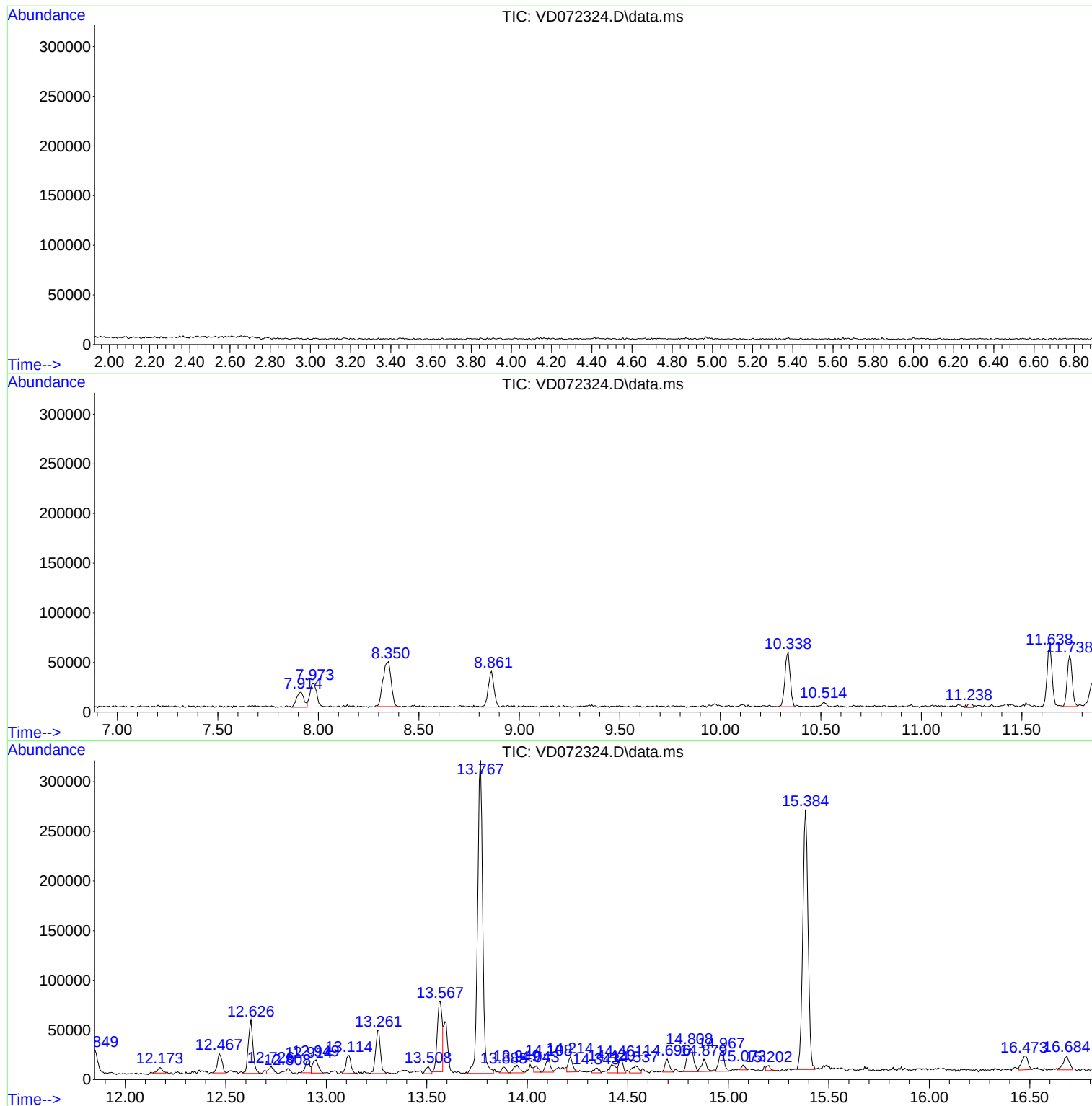
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031122S.M
Quant Title : SW846 8260

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



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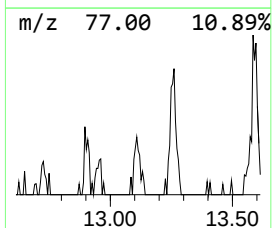
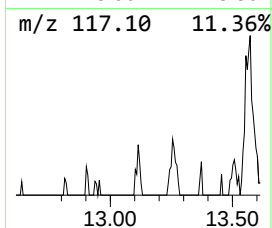
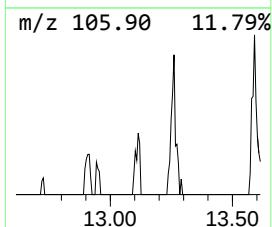
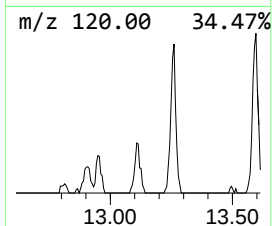
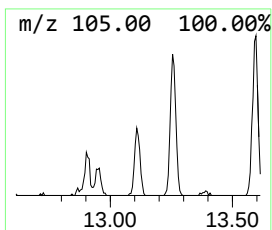
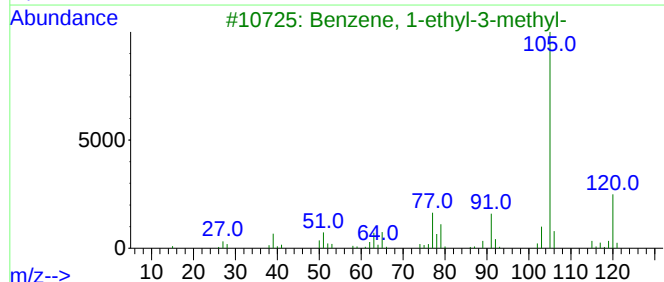
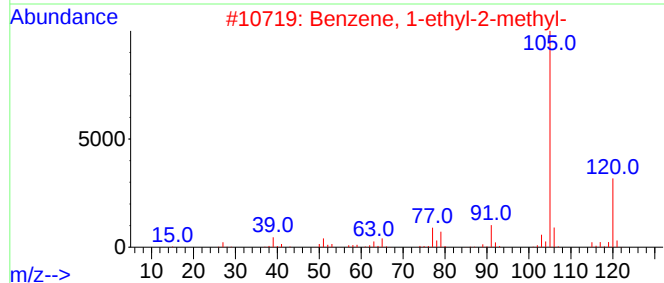
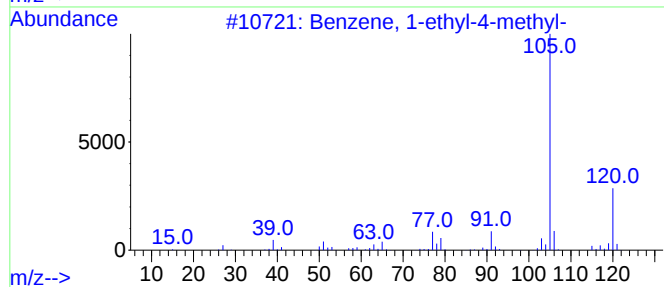
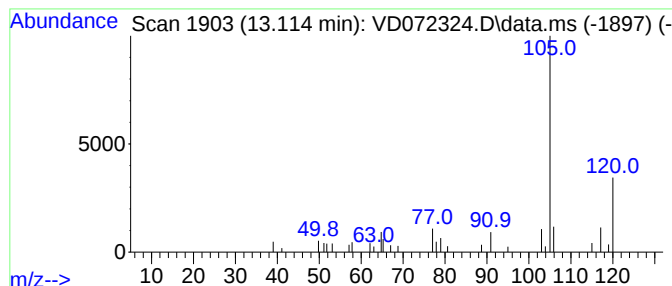
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031122S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.114	12.79 ug/l	31246	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	80
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	80
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	72
4			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	53
5			Disulfide, bis(p-methylbenzyl)	274	C16H18S2	020193-94-6	38



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ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
G-4-7-17-18

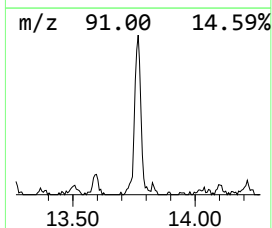
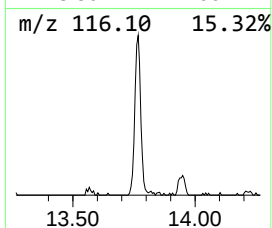
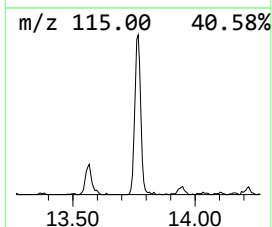
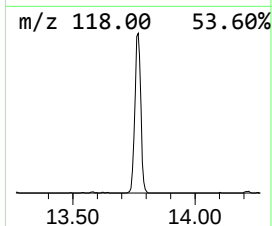
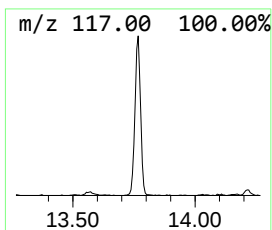
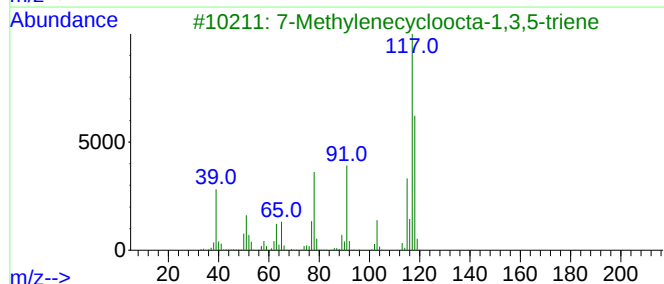
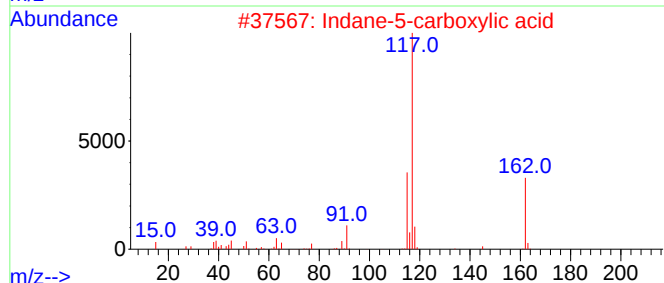
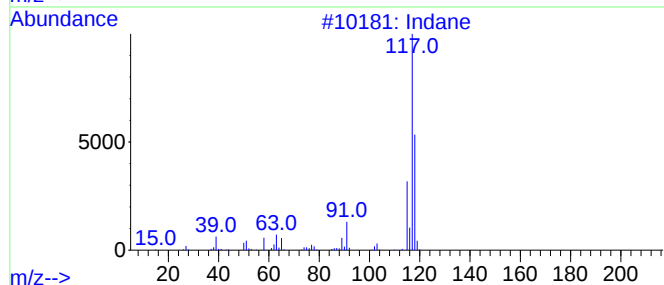
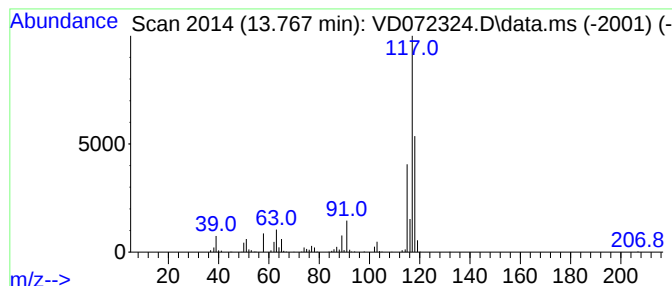
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031122S.M
Quant Title : SW846 8260

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

Peak Number 2 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.767	224.29 ug/l	547897	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	94
2			Indane-5-carboxylic acid	162	C10H10O2	065898-38-6	64
3			7-Methylenecycloocta-1,3,5-triene	118	C9H10	002570-13-0	64
4			3-Bromo-1-phenyl-1-propene	196	C9H9Br	004392-24-9	59
5			Bicyclo[4.2.0]octa-1,3,5-triene,...	118	C9H10	055337-80-9	58



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Misc : 5.06G/5.00ml/MSVOA_D/SOIL
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
G-4-7-17-18

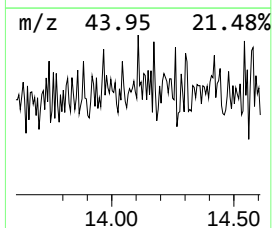
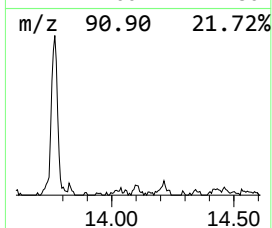
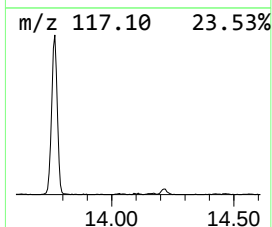
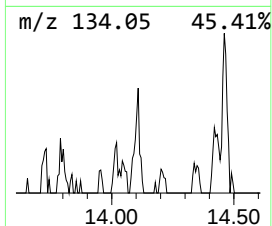
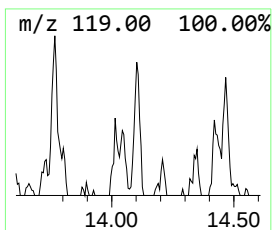
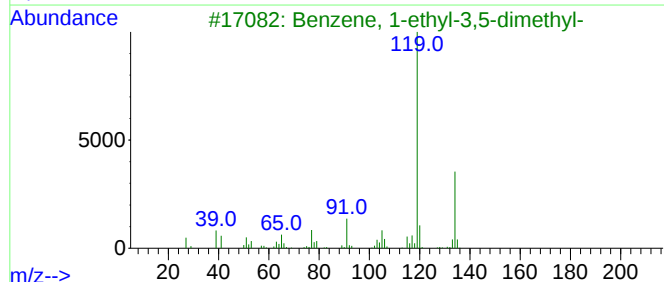
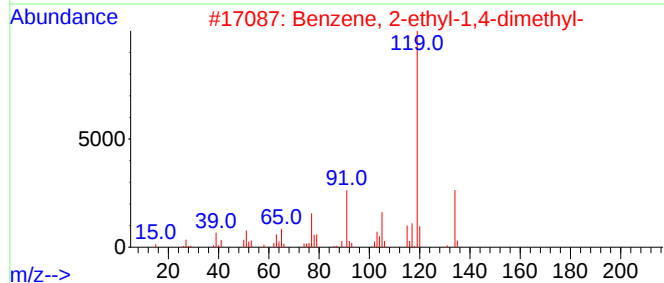
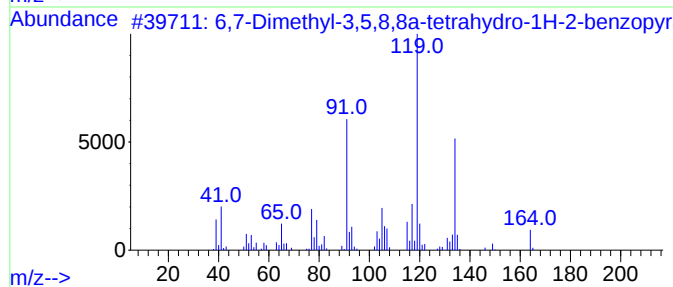
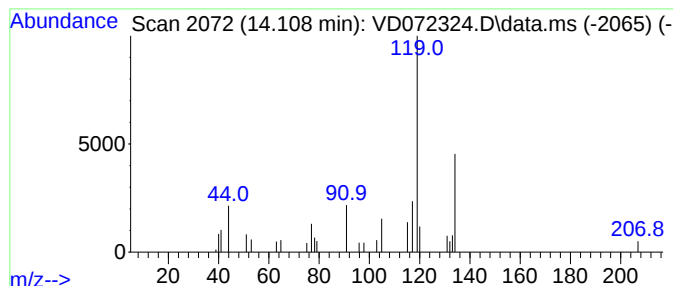
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031122S.M
Quant Title : SW846 8260

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

Peak Number 3 6,7-Dimethyl-3,5,8,8a-tetra... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.108	8.92 ug/l	21792	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6,7-Dimethyl-3,5,8,8a-tetrahydro...	164	C11H16O	110028-10-9	72
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	64
3			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	58
4			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	58
5			o-Cymene	134	C10H14	000527-84-4	58



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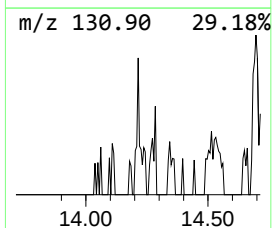
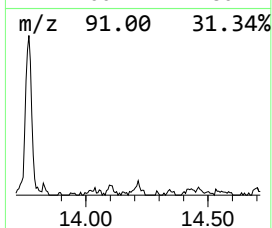
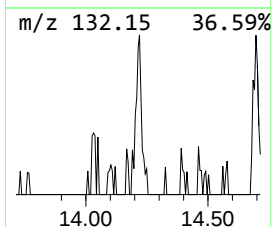
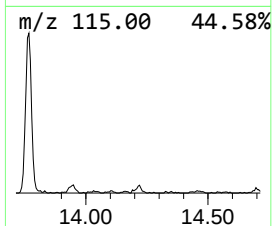
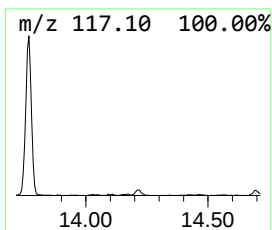
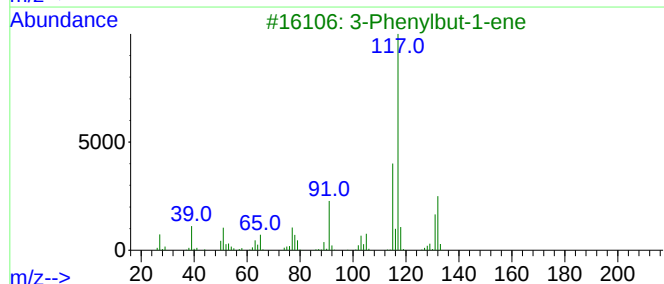
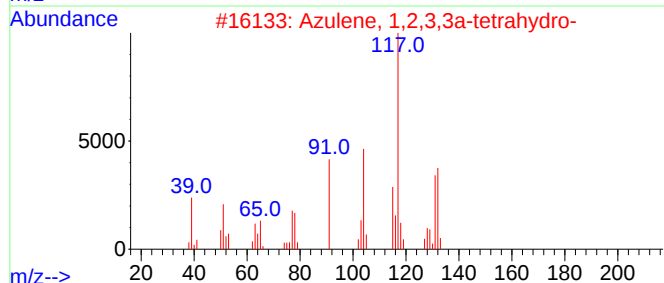
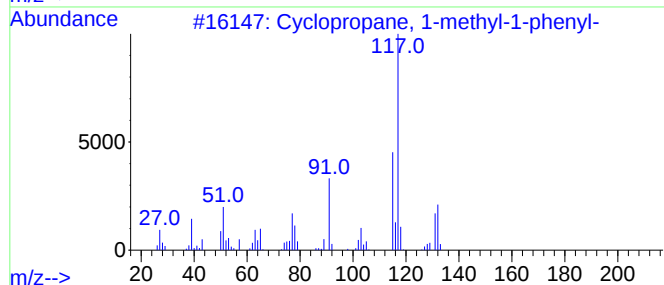
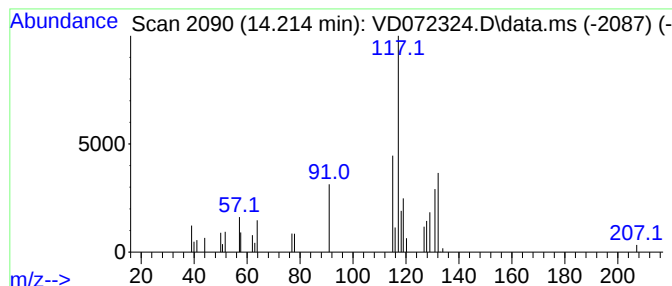
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031122S.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Cyclopropane, 1-methyl-1-ph... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.214	9.14 ug/l	22339	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropane, 1-methyl-1-phenyl-	132	C10H12	002214-14-4	59
2			Azulene, 1,2,3,3a-tetrahydro-	132	C10H12	033877-87-1	58
3			3-Phenylbut-1-ene	132	C10H12	000934-10-1	53
4			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	53
5			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	53



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ALS Vial : 46 Sample Multiplier: 1

Instrument :
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ClientSampleId :
G-4-7-17-18

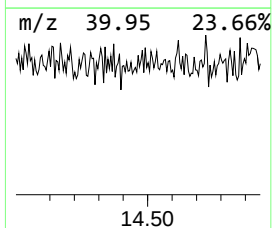
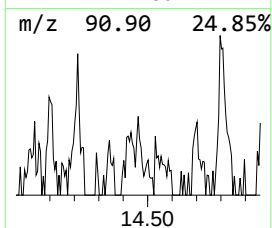
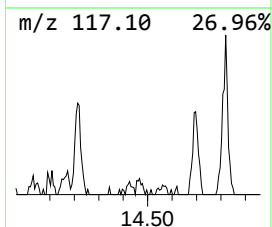
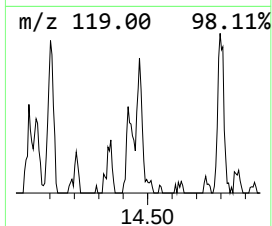
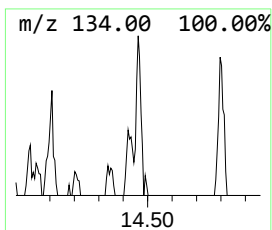
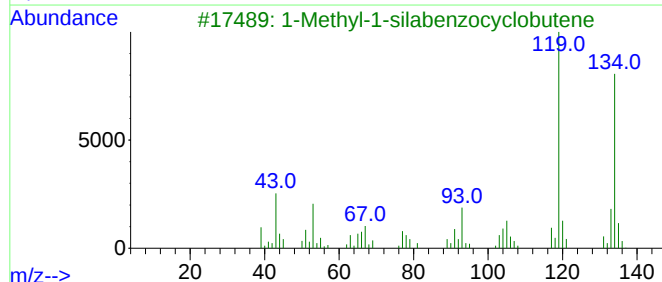
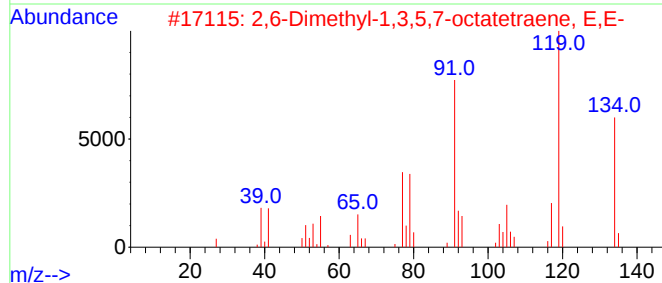
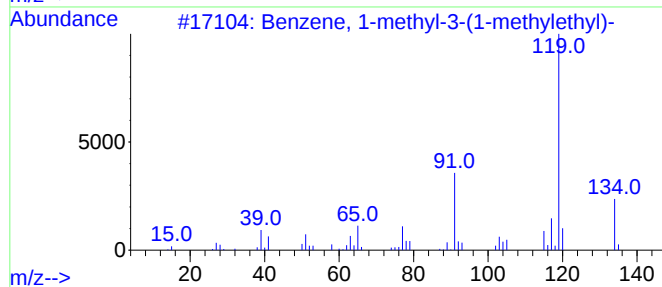
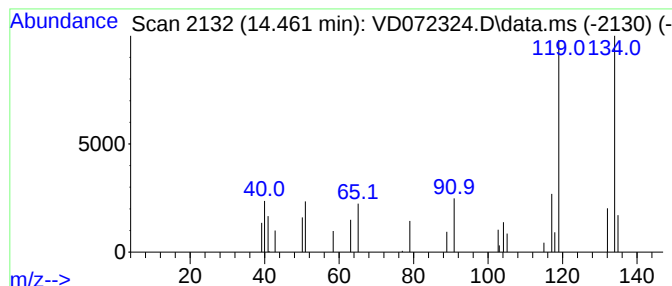
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-methyl-3-(1-meth... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.461	8.17 ug/l	19963	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	59
2			2,6-Dimethyl-1,3,5,7-octatetraen...	134	C10H14	000460-01-5	53
3			1-Methyl-1-silabenzocyclobutene	134	C8H10Si	160841-06-5	53
4			1,3,8-p-Menthatriene	134	C10H14	018368-95-1	53
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	52



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD031822\
Data File : VD072324.D
Acq On : 19 Mar 2022 08:23
Operator : VA/SY
Sample : N1972-02
Misc : 5.06G/5.00ml/MSVOA_D/SOIL
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
G-4-7-17-18

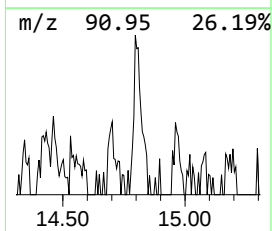
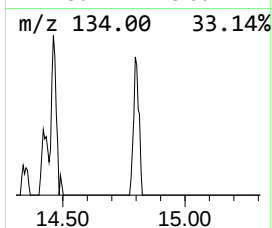
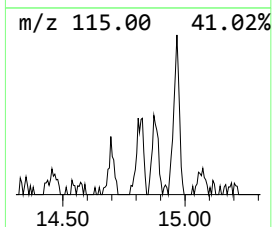
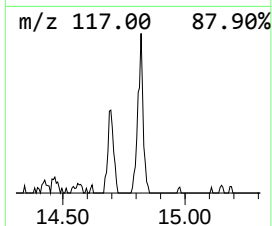
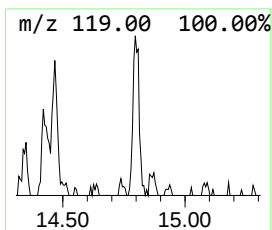
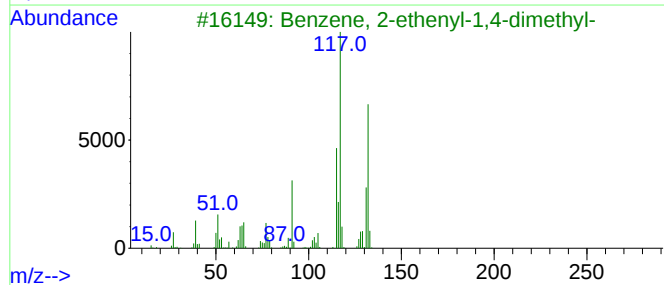
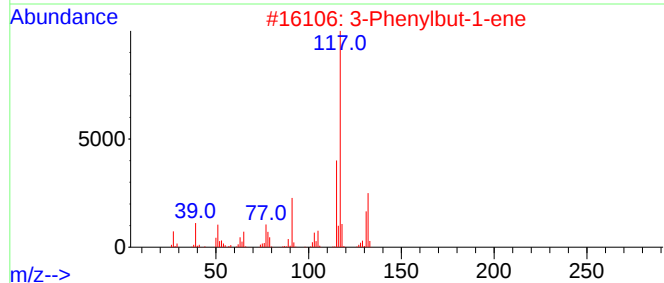
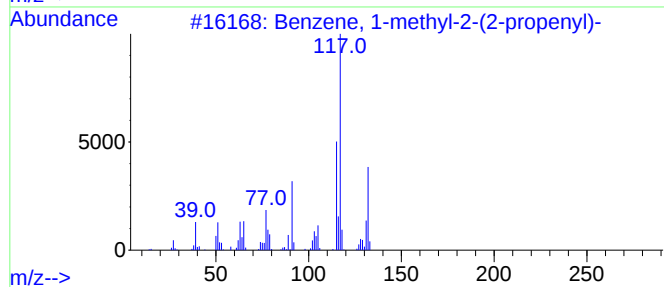
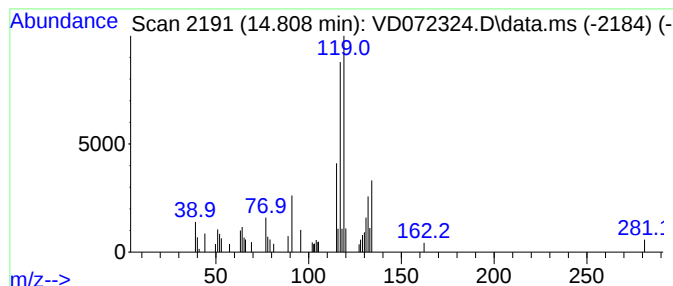
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-methyl-2-(2-prop... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.808	22.20 ug/l	54225	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	86
2			3-Phenylbut-1-ene	132	C10H12	000934-10-1	50
3			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	50
4			(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	50
5			Benzene, 2-butenyl-	132	C10H12	001560-06-1	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD031822\
Data File : VD072324.D
Acq On : 19 Mar 2022 08:23
Operator : VA/SY
Sample : N1972-02
Misc : 5.06G/5.00ml/MSVOA_D/SOIL
ALS Vial : 46 Sample Multiplier: 1

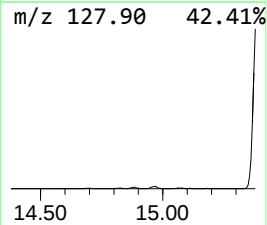
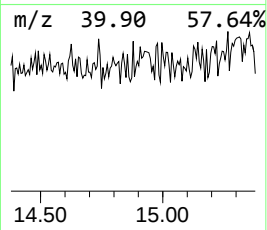
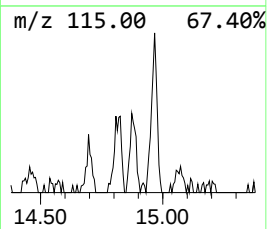
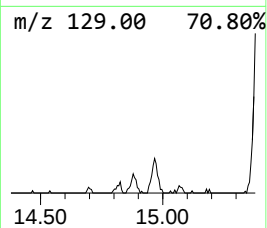
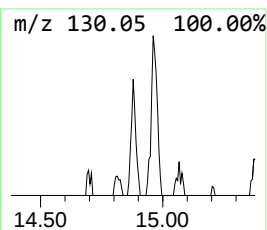
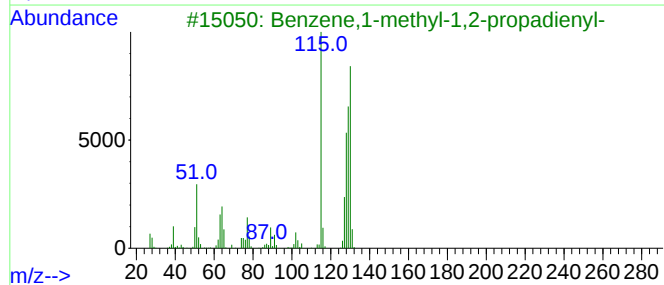
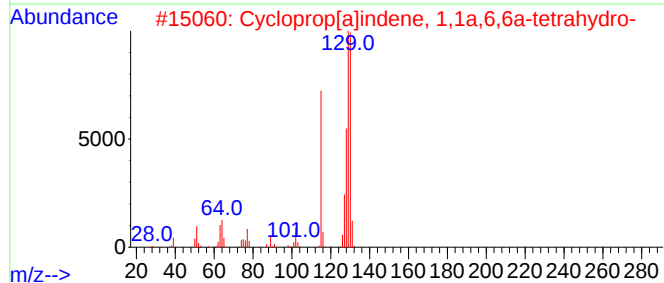
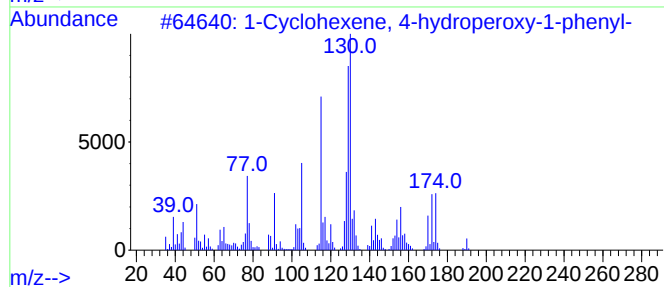
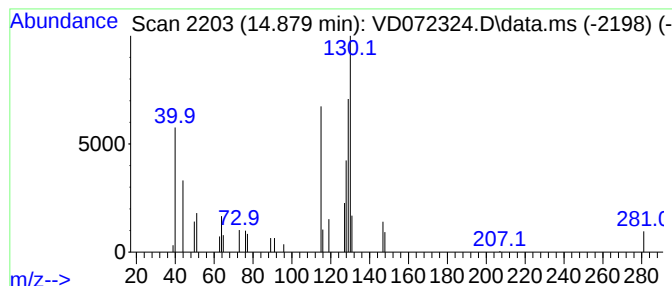
Instrument :
MSVOA_D
ClientSampleId :
G-4-7-17-18

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031122S.M
Quant Title : SW846 8260

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

Peak Number 7 1-Cyclohexene, 4-hydroperox... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.879	8.63 ug/l	21079	1,4-Dichlorobenzene-d4	13.567	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Cyclohexene, 4-hydroperoxy-1-p...	190	C12H14O2	111005-75-5	64
2	Cycloprop[a]indene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	60
3	Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	60
4	1,4-Methanonaphthalen-9-one, 1,2...	158	C11H10O	006165-88-4	59
5	Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD031822\
Data File : VD072324.D
Acq On : 19 Mar 2022 08:23
Operator : VA/SY
Sample : N1972-02
Misc : 5.06G/5.00ml/MSVOA_D/SOIL
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
G-4-7-17-18

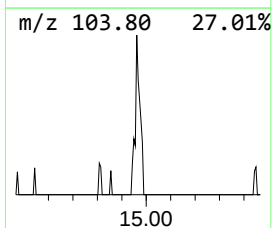
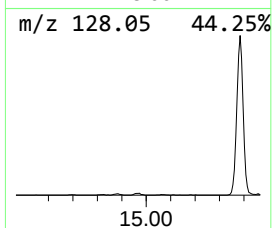
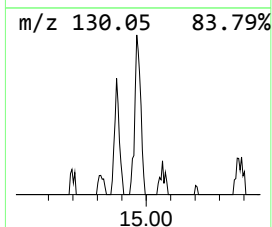
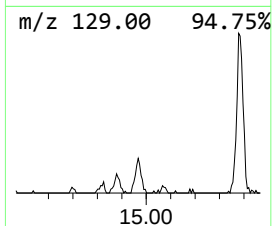
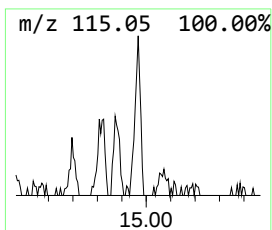
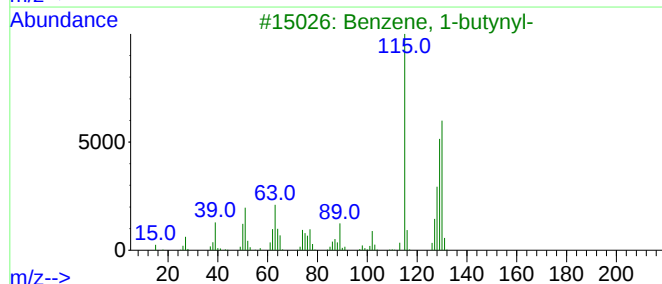
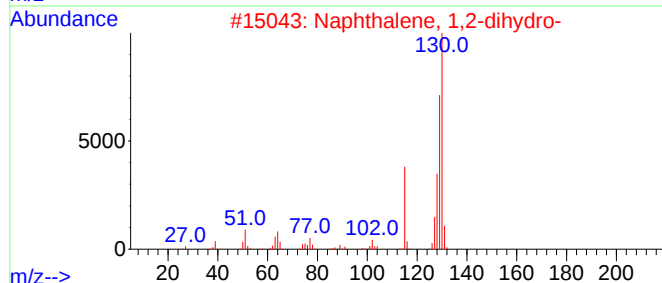
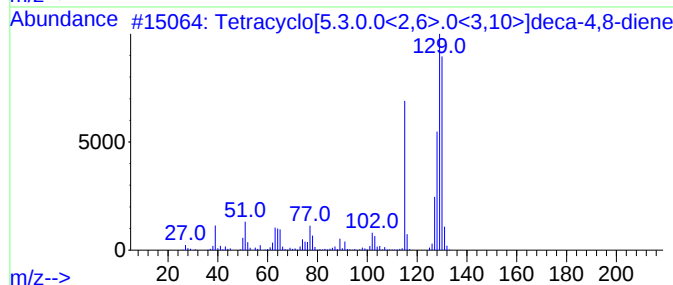
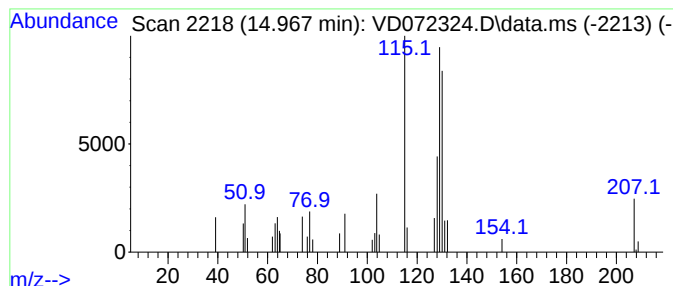
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Quant Title : SW846 8260

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

Peak Number 8 Tetracyclo[5.3.0.0<2,6>.0<3... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.967	15.18 ug/l	37076	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracyclo[5.3.0.0<2,6>.0<3,10>]...	130	C10H10	034324-40-8	86
2			Naphthalene, 1,2-dihydro-	130	C10H10	000447-53-0	76
3			Benzene, 1-butynyl-	130	C10H10	000622-76-4	70
4			Cycloprop[a]indene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	62
5			Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD031822\
Data File : VD072324.D
Acq On : 19 Mar 2022 08:23
Operator : VA/SY
Sample : N1972-02
Misc : 5.06G/5.00ml/MSVOA_D/SOIL
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
G-4-7-17-18

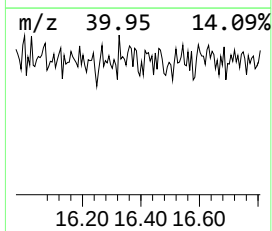
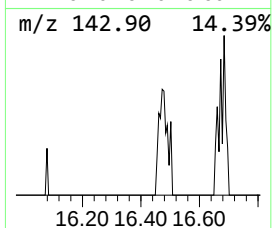
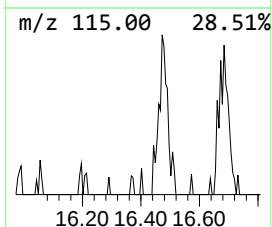
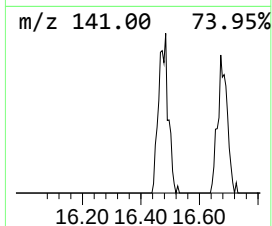
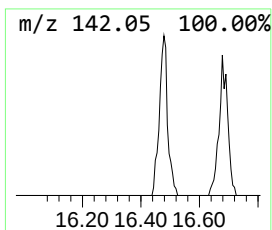
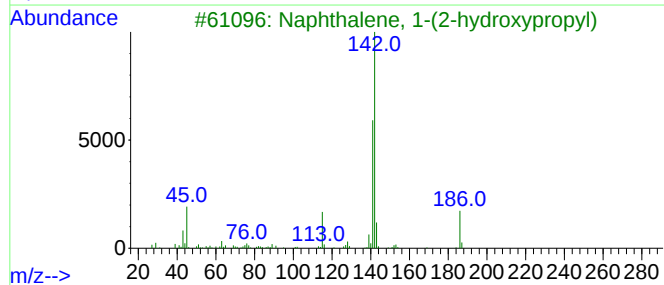
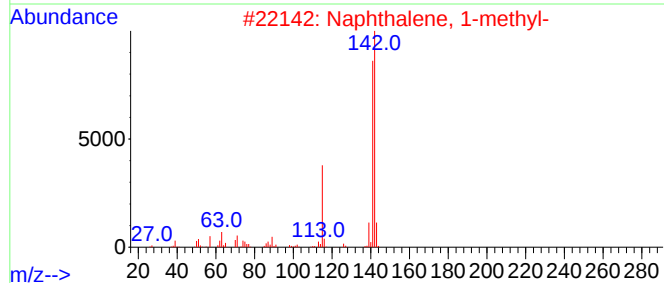
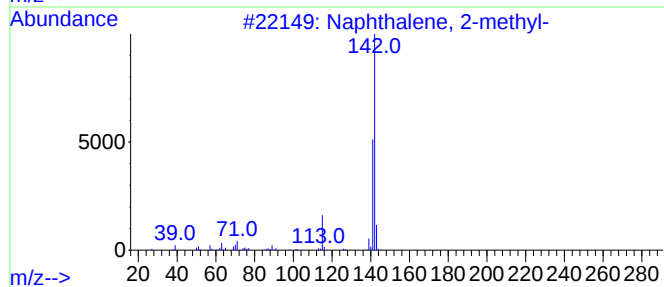
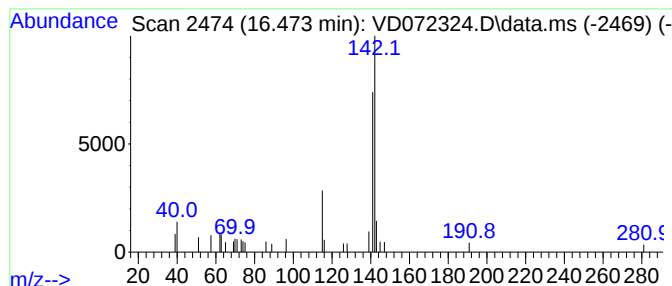
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-(2-hydroxypr... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.473	12.26 ug/l	29947	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	83
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	80
3			Naphthalene, 1-(2-hydroxypropyl)	186	C13H14O	027653-13-0	78
4			Benzocycloheptatriene	142	C11H10	000264-09-5	72
5			Spiro(9-methylenetricyclo[6.2.1]....	184	C13H12O	033929-47-4	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD031822\
Data File : VD072324.D
Acq On : 19 Mar 2022 08:23
Operator : VA/SY
Sample : N1972-02
Misc : 5.06G/5.00ml/MSVOA_D/SOIL
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
G-4-7-17-18

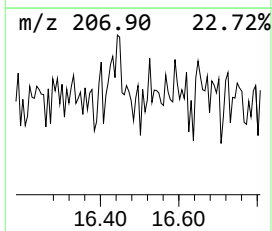
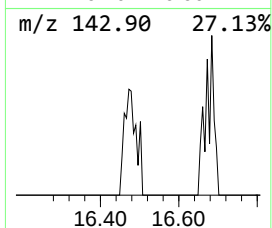
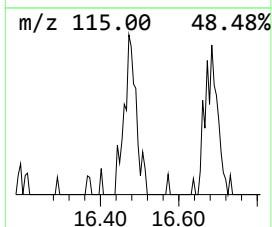
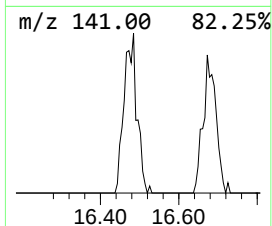
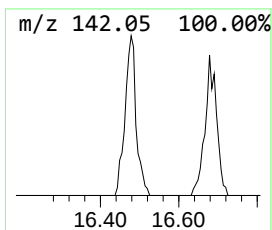
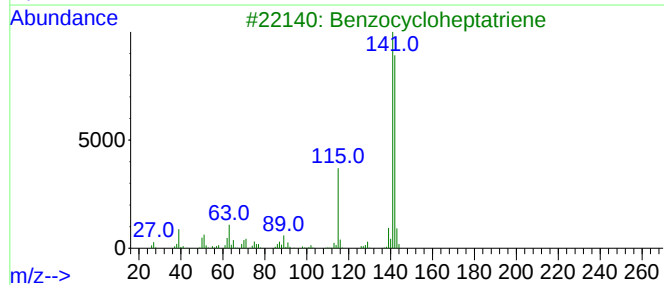
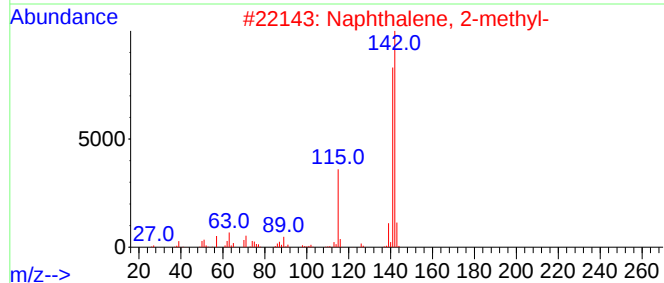
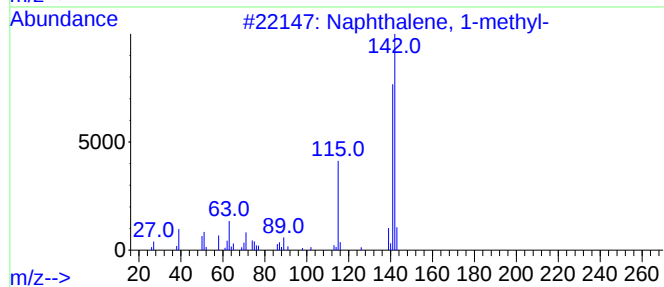
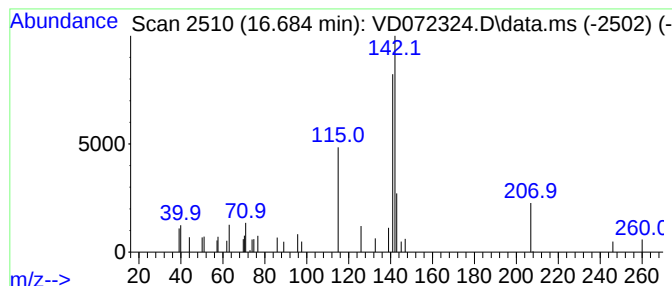
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Quant Title : SW846 8260

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

Peak Number 10 Benzocycloheptatriene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.684	12.21 ug/l	29828	1,4-Dichlorobenzene-d4	13.567

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	72
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	64
3			Benzocycloheptatriene	142	C11H10	000264-09-5	58
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	58
5			3,4-Difluorobenzaldehyde	142	C7H4F2O	034036-07-2	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD031822\
Data File : VD072324.D
Acq On : 19 Mar 2022 08:23
Operator : VA/SY
Sample : N1972-02
Misc : 5.06G/5.00ml/MSVOA_D/SOIL
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
G-4-7-17-18

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031122S.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
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Indane	13.767	224.3	ug/l	547897	4	13.567	122140	50.0
6,7-Dimethyl-3,...	14.108	8.9	ug/l	21792	4	13.567	122140	50.0
Cyclopropane, 1...	14.214	9.1	ug/l	22339	4	13.567	122140	50.0
Benzene, 1-meth...	14.461	8.2	ug/l	19963	4	13.567	122140	50.0
Benzene, 1-meth...	14.808	22.2	ug/l	54225	4	13.567	122140	50.0
1-Cyclohexene, ...	14.879	8.6	ug/l	21079	4	13.567	122140	50.0
Tetracyclo[5.3....	14.967	15.2	ug/l	37076	4	13.567	122140	50.0
Naphthalene, 1-...	16.473	12.3	ug/l	29947	4	13.567	122140	50.0
Benzocyclohepta...	16.684	12.2	ug/l	29828	4	13.567	122140	50.0