

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041824\
 Data File : VY017982.D
 Acq On : 18 Apr 2024 17:38
 Operator : SY/MD
 Sample : P2178-01
 Misc : 4.17g/5.0mL/MSVOA_Y/SOIL/A
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 TOPSOIL-LKWD-DENALI-4

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_Y\methods\82Y041624S.M
 Title : SW846 8260

Signal : TIC: VY017982.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.387	87	93	101	rBV	36833	66198	1.02%	0.175%
2	3.954	341	350	360	rBV2	26385	74671	1.15%	0.198%
3	7.722	958	968	974	rBV	253790	612448	9.42%	1.622%
4	7.795	974	980	997	rVB	389090	884197	13.61%	2.342%
5	8.148	1025	1038	1050	rBV	210510	477622	7.35%	1.265%
6	8.697	1119	1128	1140	rBV	651524	1279339	19.69%	3.389%
7	9.270	1214	1222	1231	rBV	38895	74853	1.15%	0.198%
8	10.185	1364	1372	1379	rBV	1012271	1789559	27.54%	4.740%
9	10.941	1490	1496	1512	rVB	275084	480292	7.39%	1.272%
10	11.495	1580	1587	1599	rBV	886513	1446768	22.26%	3.832%
11	11.715	1611	1623	1628	rBV5	26641	89263	1.37%	0.236%
12	11.898	1637	1653	1655	rBV2	129767	496363	7.64%	1.315%
13	12.416	1714	1738	1746	rBV2	754250	4685129	72.09%	12.410%
14	12.489	1746	1750	1771	rVB3	1164298	2925630	45.02%	7.750%
15	12.971	1808	1829	1840	rVV	1366547	6498891	100.00%	17.215%
16	13.111	1840	1852	1870	rVB2	646386	3375766	51.94%	8.942%
17	13.434	1898	1905	1911	rBV	607190	936466	14.41%	2.481%
18	13.580	1916	1929	1935	rBV4	146561	611848	9.41%	1.621%
19	13.666	1935	1943	1958	rVB	358606	1372150	21.11%	3.635%
20	14.038	1996	2004	2028	rVB2	117841	461046	7.09%	1.221%
21	14.306	2036	2048	2064	rVB2	116565	409989	6.31%	1.086%
22	14.550	2079	2088	2096	rVB4	31216	100897	1.55%	0.267%
23	14.739	2106	2119	2132	rBV	1638222	5041038	77.57%	13.353%
24	14.867	2132	2140	2149	rVV2	89884	305979	4.71%	0.811%
25	14.958	2151	2155	2163	rVB5	41917	92614	1.43%	0.245%
26	15.056	2163	2171	2176	rBV3	32161	89572	1.38%	0.237%
27	15.129	2176	2183	2193	rVV2	85269	264368	4.07%	0.700%
28	15.239	2193	2201	2202	rVV4	28011	73095	1.12%	0.194%
29	15.293	2203	2210	2221	rVV4	82291	315505	4.85%	0.836%
30	15.440	2221	2234	2254	rVB	438050	1578876	24.29%	4.182%
31	15.745	2273	2284	2296	rVB3	50553	189612	2.92%	0.502%
32	15.915	2303	2312	2325	rVB3	35217	114568	1.76%	0.303%
33	16.238	2355	2365	2369	rVV7	28272	94054	1.45%	0.249%
34	16.299	2369	2375	2382	rVB	108453	228593	3.52%	0.606%
35	16.494	2399	2407	2416	rBV	98987	214096	3.29%	0.567%

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Peak separation: 5

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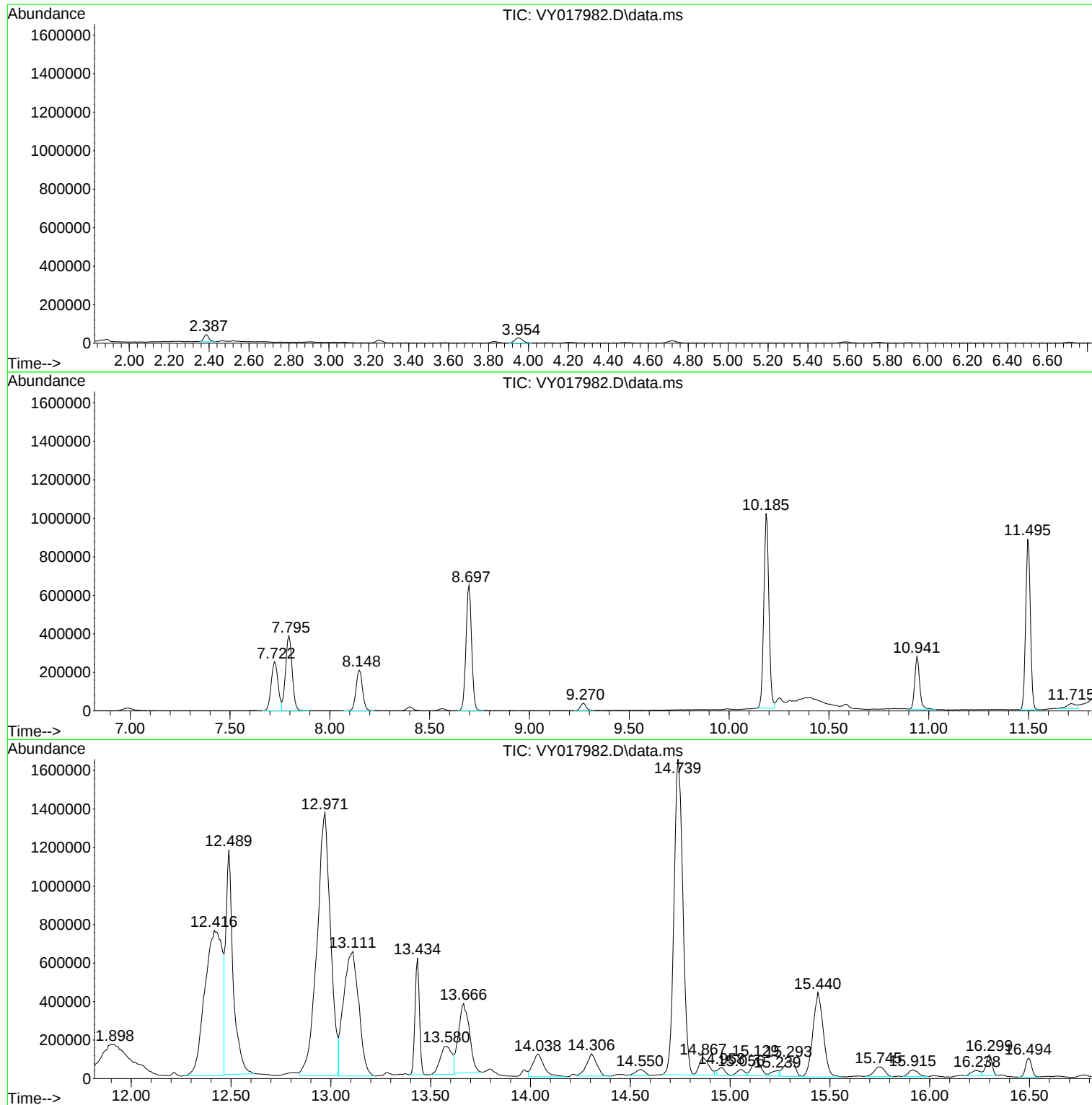
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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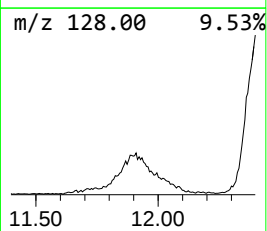
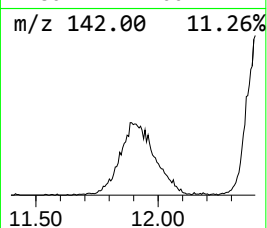
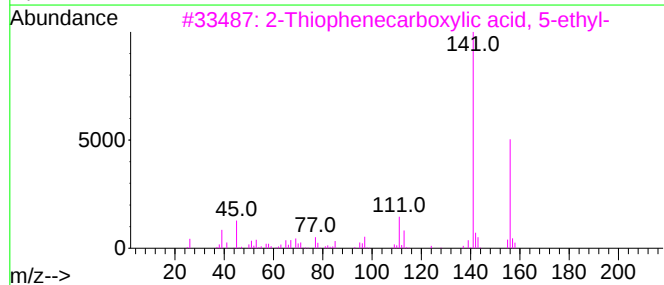
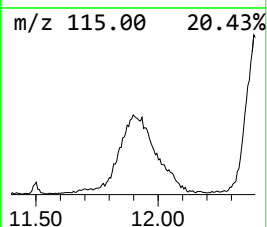
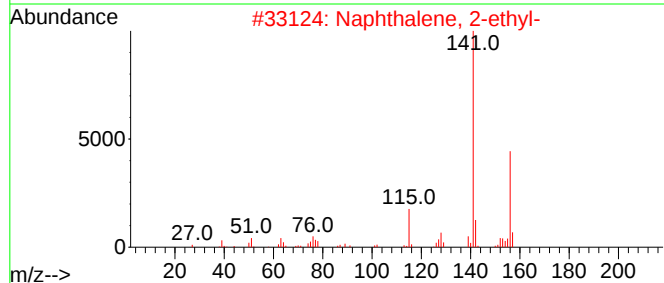
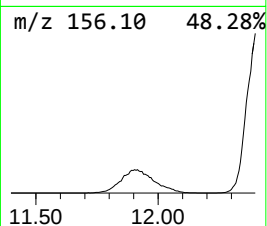
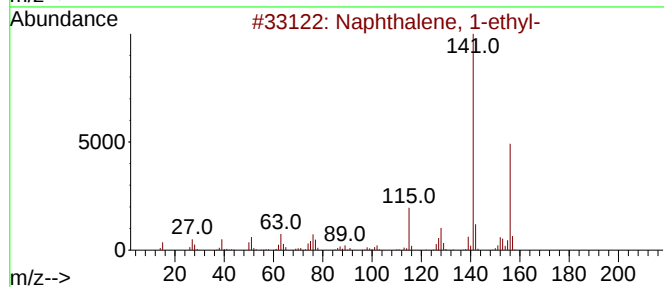
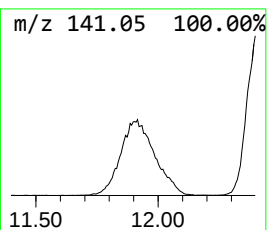
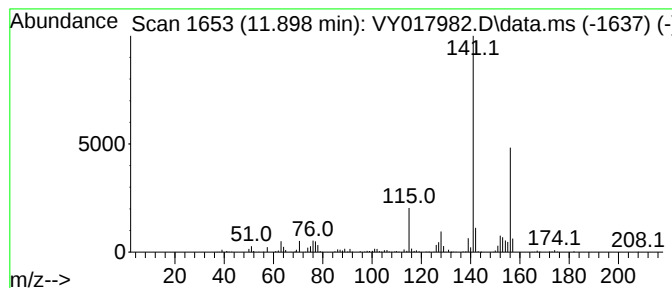
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041624S.M
Quant Title : SW846 8260

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

Peak Number 1 Naphthalene, 1-ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.898	17.15 ug/l	496363	Chlorobenzene-d5	11.495

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	97
2			Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	97
3			2-Thiophenecarboxylic acid, 5-et...	156	C7H8O2S	023229-72-3	59
4			2,4-Difluoromesitylene	156	C9H10F2	000392-61-0	59
5			5-Acetyl-2-amino-4-methylthiazole	156	C6H8N2OS	030748-47-1	53



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Instrument :
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ClientSampleId :
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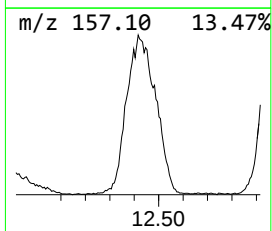
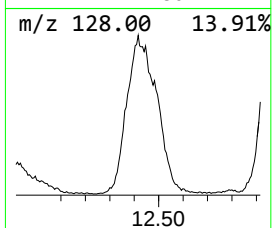
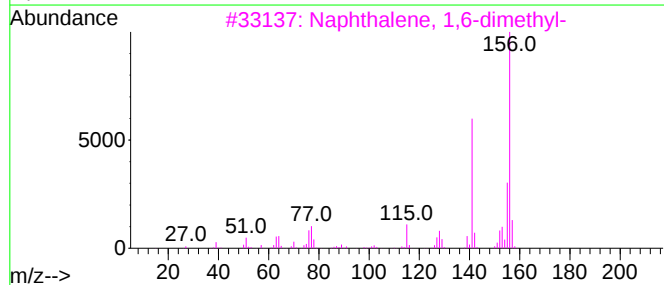
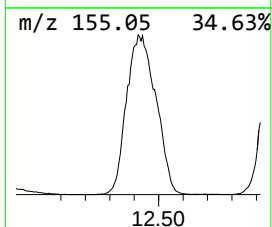
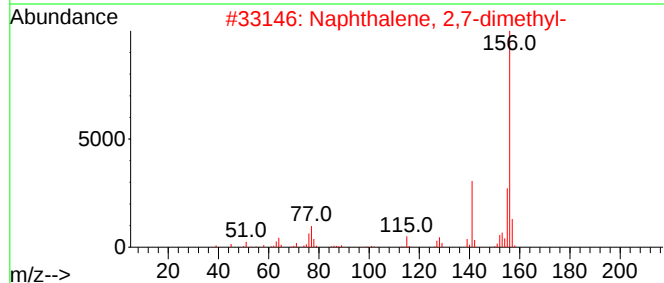
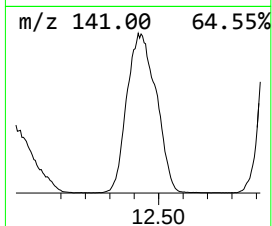
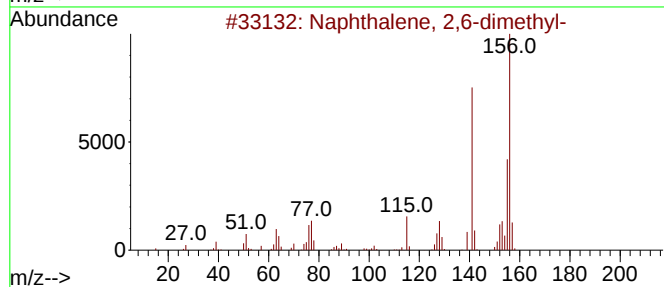
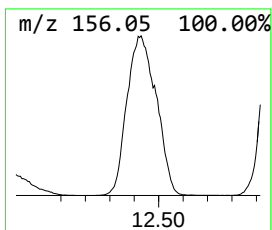
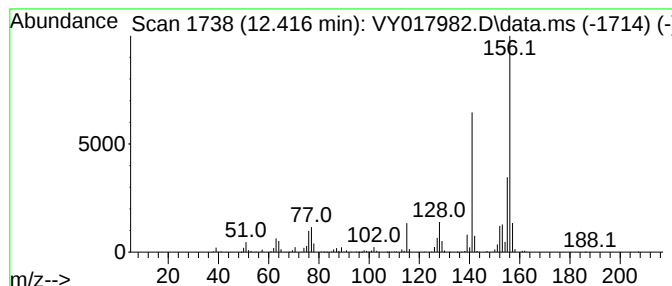
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041624S.M
Quant Title : SW846 8260

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

Peak Number 2 Naphthalene, 2,6-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.416	161.92 ug/l	4685130	Chlorobenzene-d5	11.495

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	98
2			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
4			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
5			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97



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Instrument :
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ClientSampleId :
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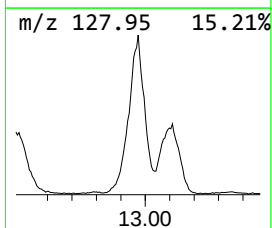
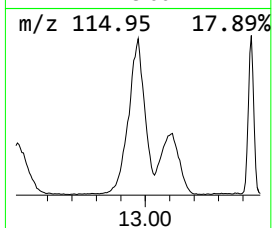
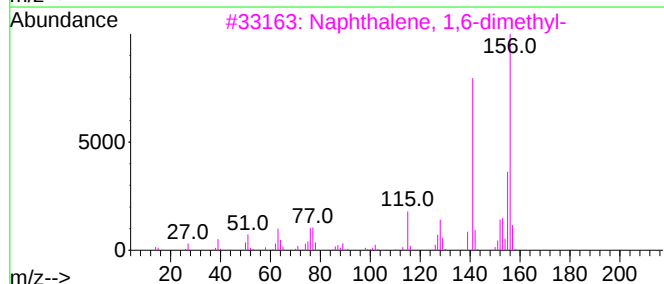
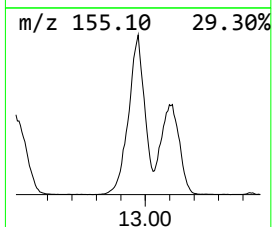
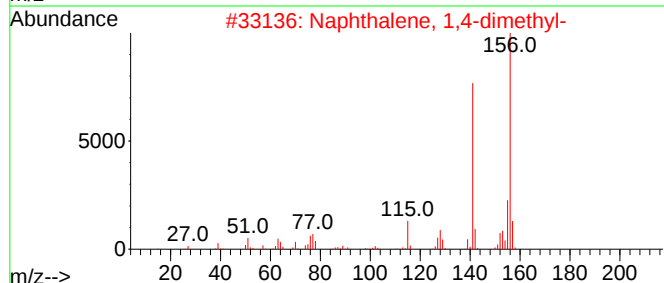
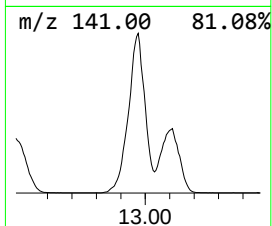
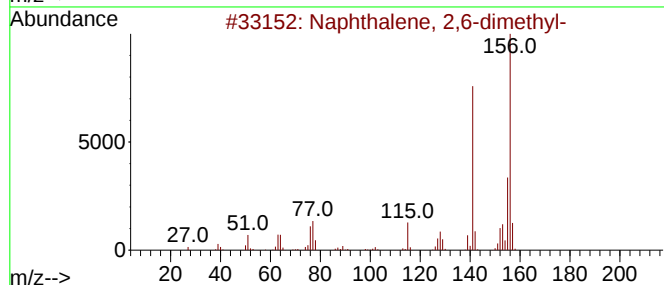
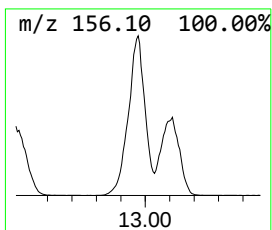
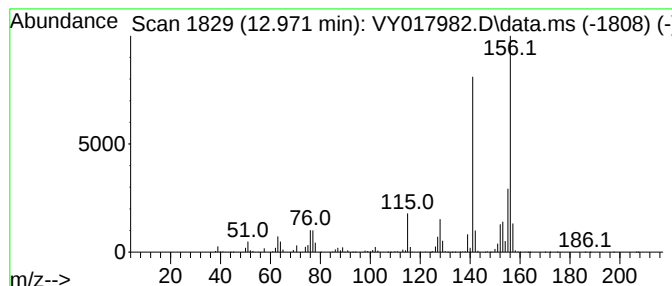
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TIC Library : C:\Database\NIST20.L
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Peak Number 3 Naphthalene, 1,4-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.971	346.99 ug/l	6498890	1,4-Dichlorobenzene-d4	13.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
2			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
5			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97



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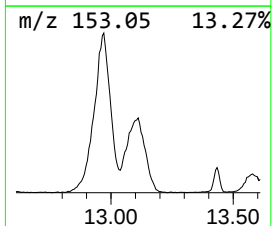
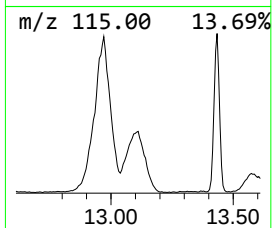
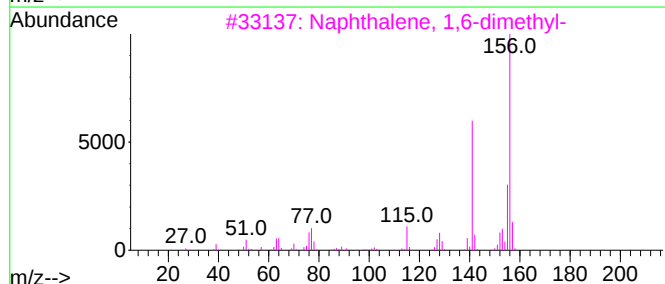
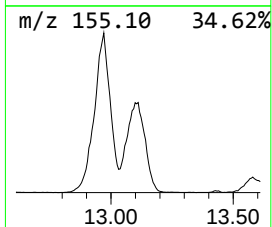
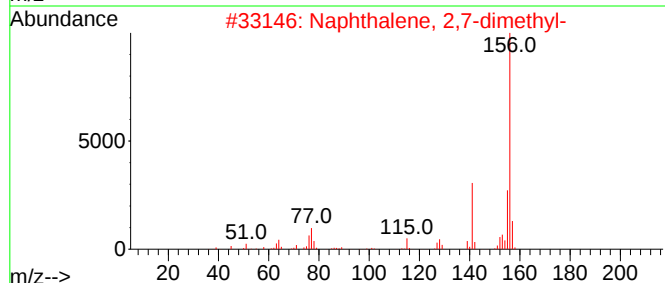
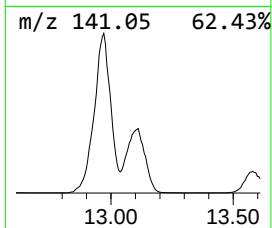
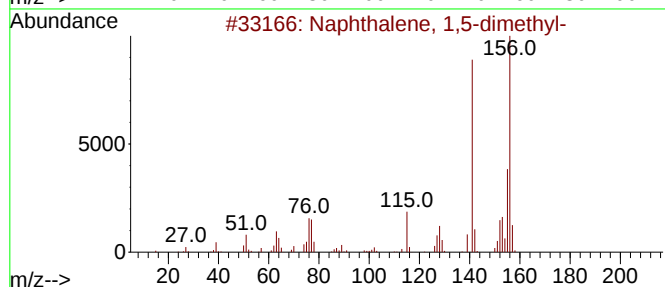
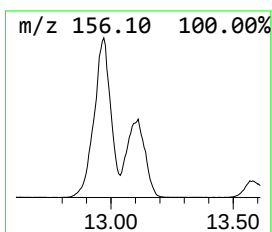
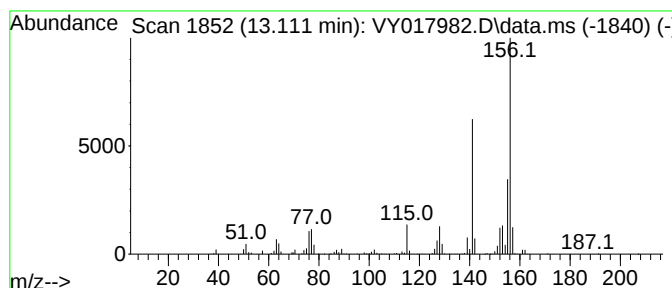
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Peak Number 4 Naphthalene, 1,5-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.111	180.24 ug/l	3375770	1,4-Dichlorobenzene-d4	13.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	98
2			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
4			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
5			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97



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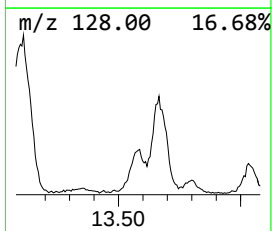
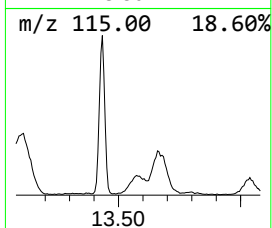
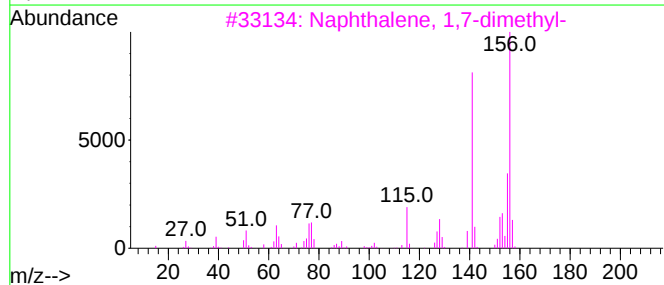
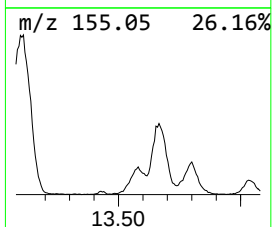
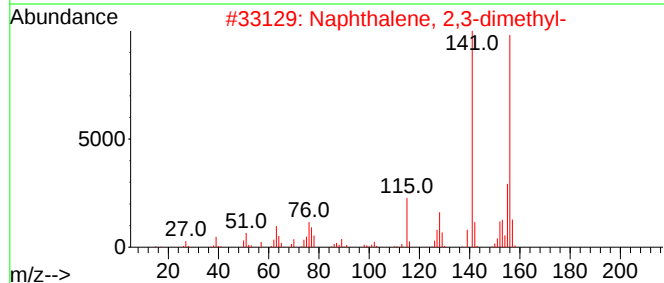
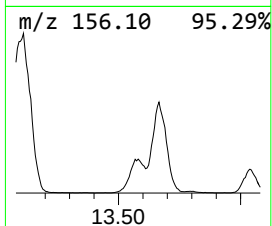
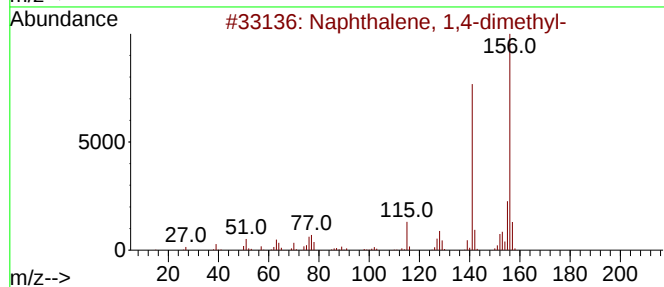
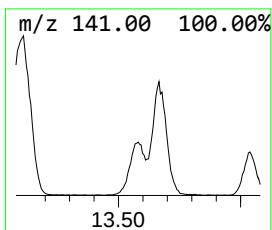
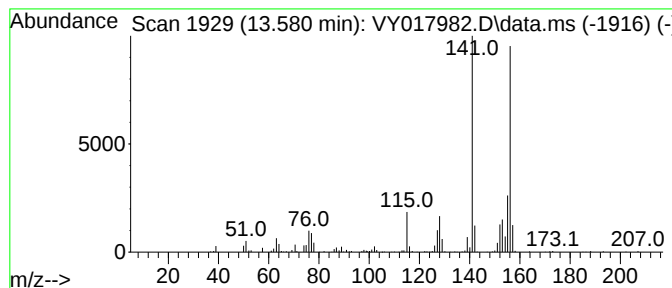
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TIC Integration Parameters: LSCINT.P

Peak Number 5 Naphthalene, 2,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.580	32.67 ug/l	611848	1,4-Dichlorobenzene-d4	13.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
3			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95
4			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
5			Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041824\
Data File : VY017982.D
Acq On : 18 Apr 2024 17:38
Operator : SY/MD
Sample : P2178-01
Misc : 4.17g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TOPSOIL-LKWD-DENALI-4

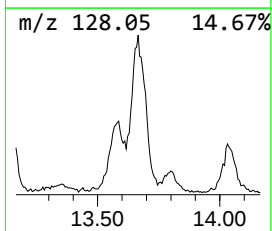
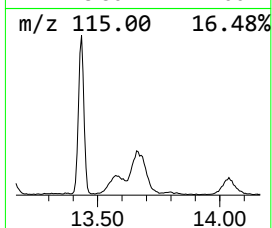
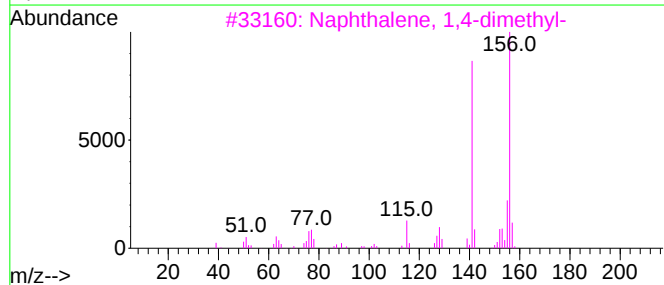
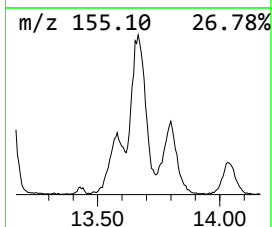
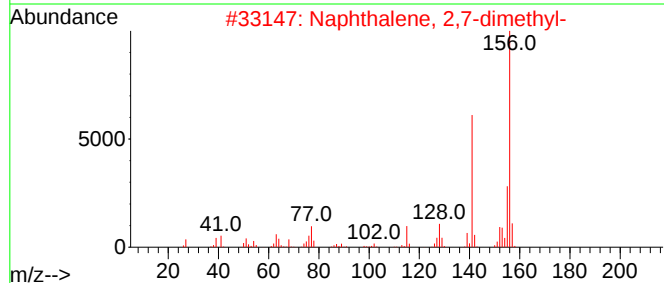
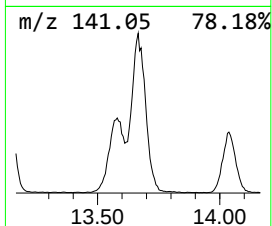
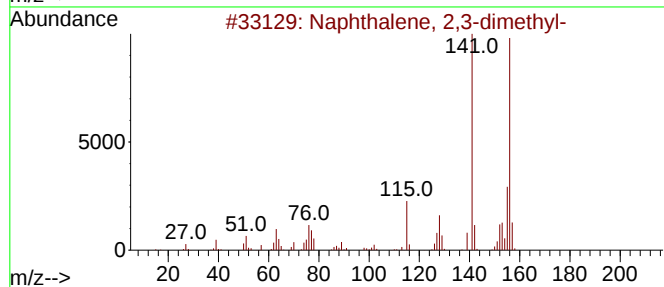
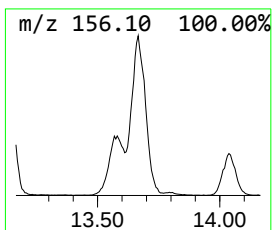
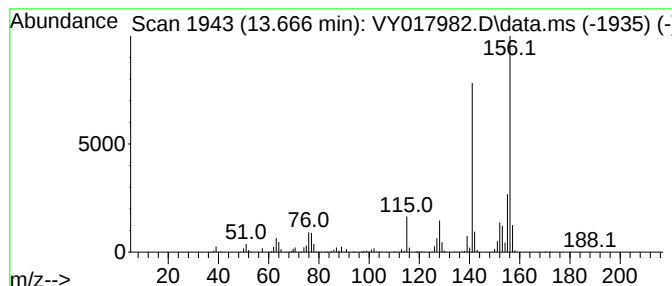
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Quant Title : SW846 8260

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

Peak Number 6 Naphthalene, 2,7-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.666	73.26 ug/l	1372150	1,4-Dichlorobenzene-d4	13.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
4			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
5			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041824\
Data File : VY017982.D
Acq On : 18 Apr 2024 17:38
Operator : SY/MD
Sample : P2178-01
Misc : 4.17g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TOPSOIL-LKWD-DENALI-4

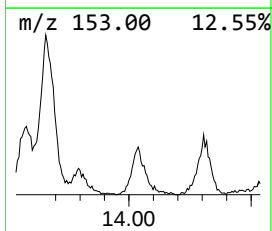
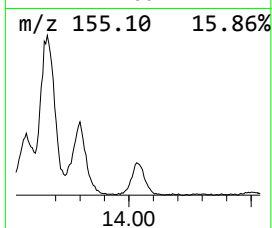
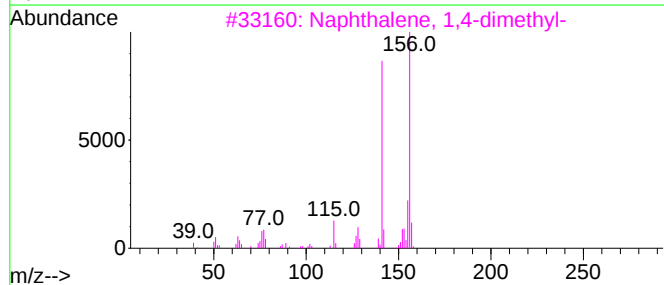
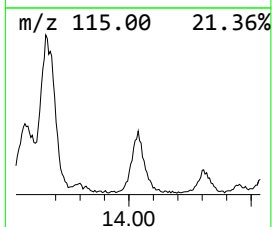
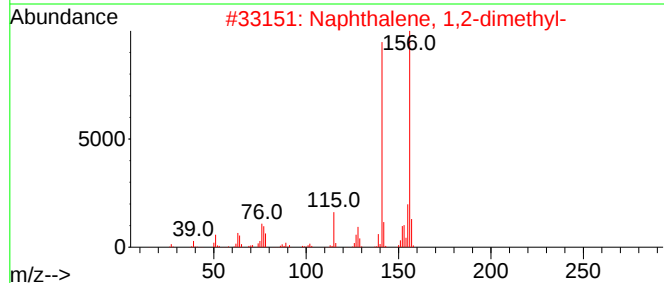
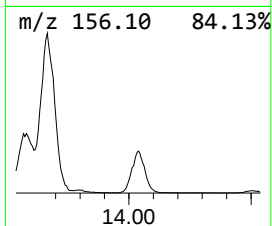
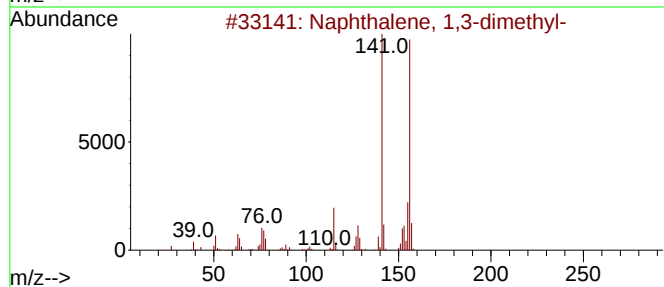
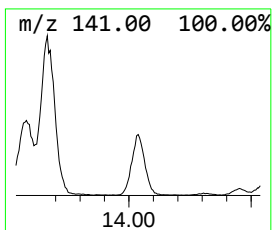
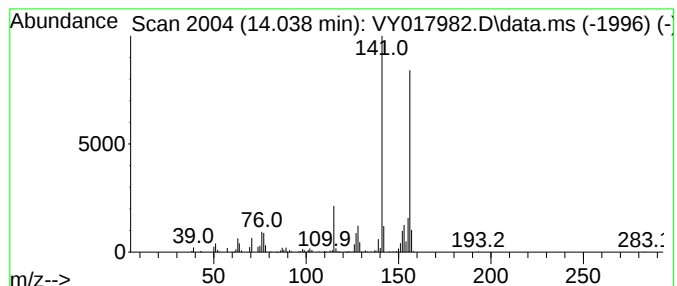
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Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 1,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.038	24.62 ug/l	461046	1,4-Dichlorobenzene-d4	13.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
2			Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	96
3			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95
5			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041824\
Data File : VY017982.D
Acq On : 18 Apr 2024 17:38
Operator : SY/MD
Sample : P2178-01
Misc : 4.17g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TOPSOIL-LKWD-DENALI-4

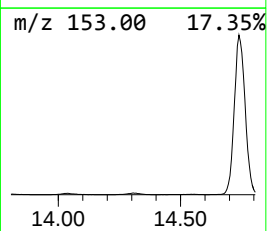
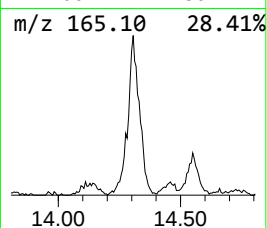
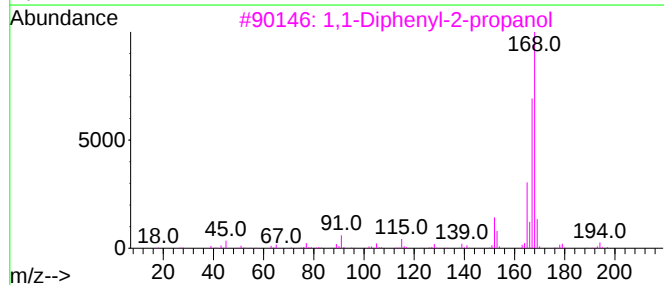
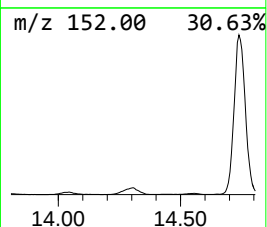
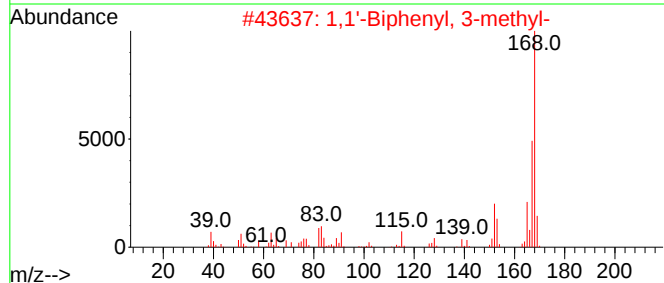
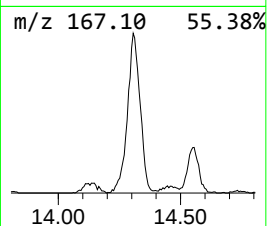
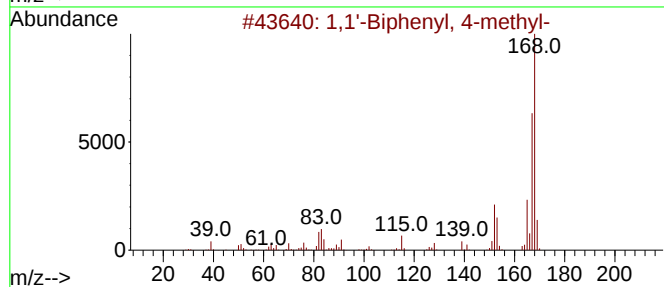
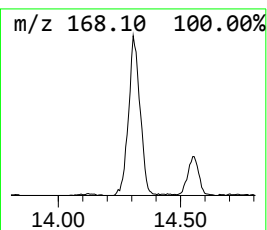
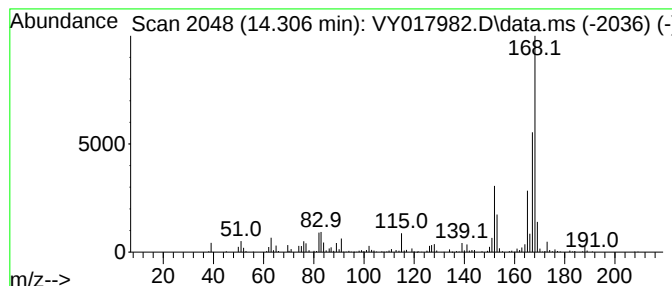
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Quant Title : SW846 8260

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

Peak Number 8 1,1'-Biphenyl, 4-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.306	21.89 ug/l	409989	1,4-Dichlorobenzene-d4	13.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,1'-Biphenyl, 4-methyl-	168	C13H12	000644-08-6	95
2			1,1'-Biphenyl, 3-methyl-	168	C13H12	000643-93-6	94
3			1,1-Diphenyl-2-propanol	212	C15H16O	029338-49-6	53
4			Nordiphenamid	225	C15H15NO	000954-21-2	52
5			1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041824\
Data File : VY017982.D
Acq On : 18 Apr 2024 17:38
Operator : SY/MD
Sample : P2178-01
Misc : 4.17g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TOPSOIL-LKWD-DENALI-4

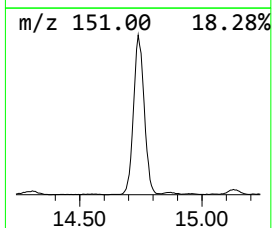
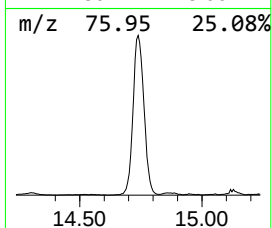
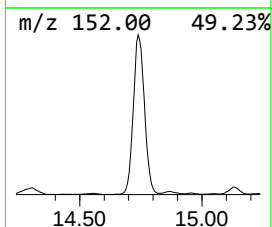
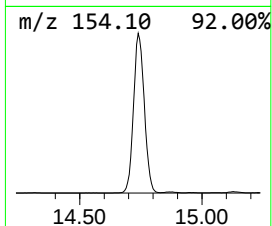
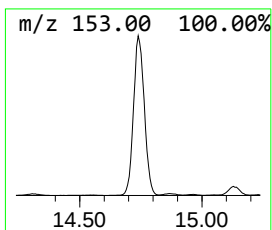
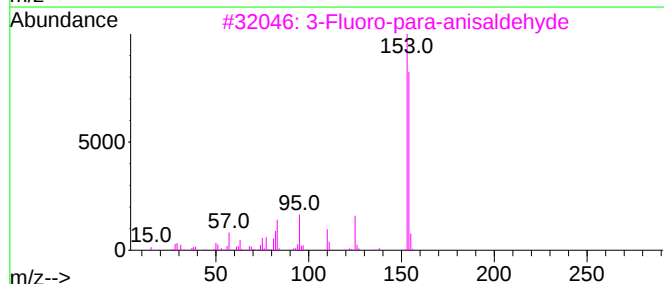
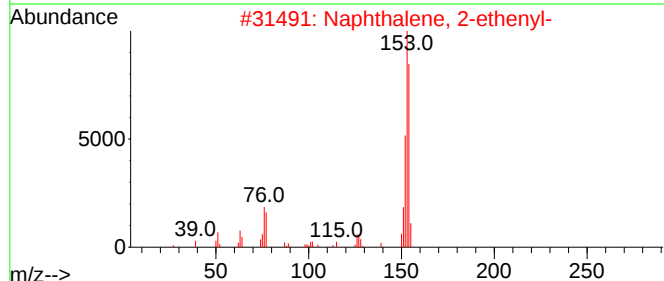
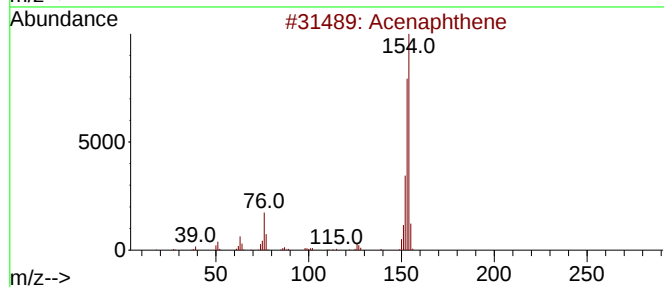
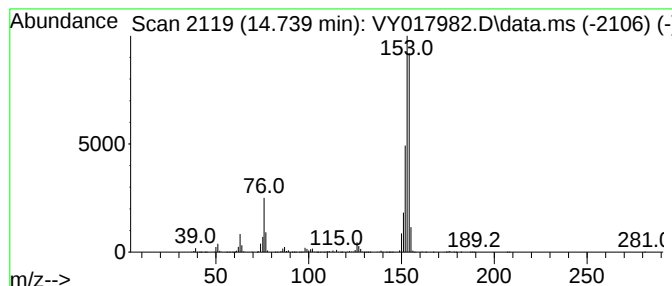
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Quant Title : SW846 8260

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

Peak Number 9 3-Fluoro-para-anisaldehyde Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.739	269.15 ug/l	5041040	1,4-Dichlorobenzene-d4	13.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acenaphthene	154	C12H10	000083-32-9	93
2			Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	59
3			3-Fluoro-para-anisaldehyde	154	C8H7FO2	000351-54-2	53
4			Biphenyl	154	C12H10	000092-52-4	43
5			1,4-Ethenonaphthalene, 1,4-dihydro-	154	C12H10	007322-47-6	42



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041824\
Data File : VY017982.D
Acq On : 18 Apr 2024 17:38
Operator : SY/MD
Sample : P2178-01
Misc : 4.17g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TOPSOIL-LKWD-DENALI-4

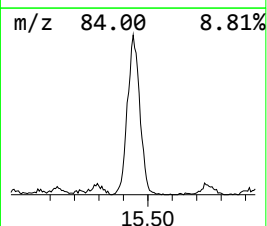
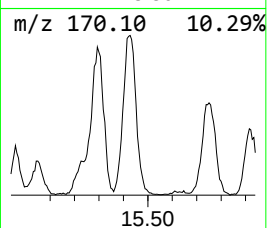
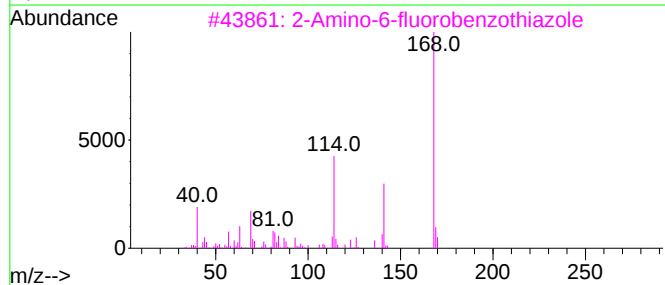
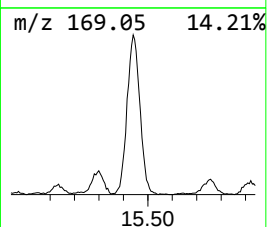
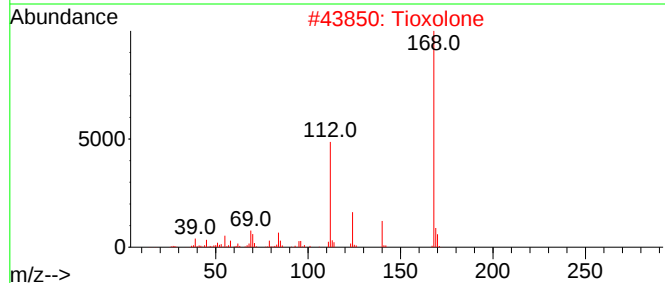
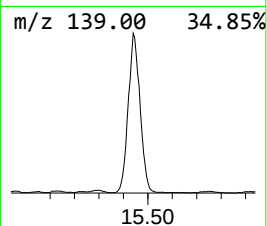
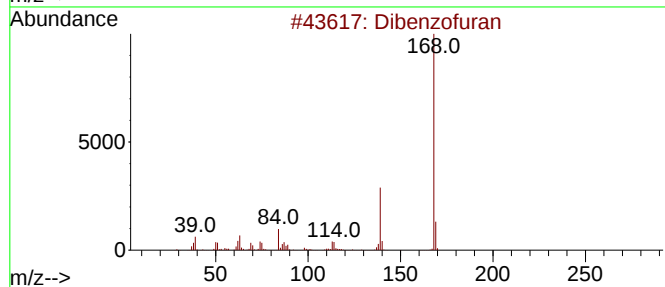
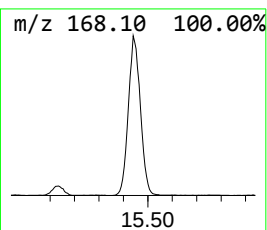
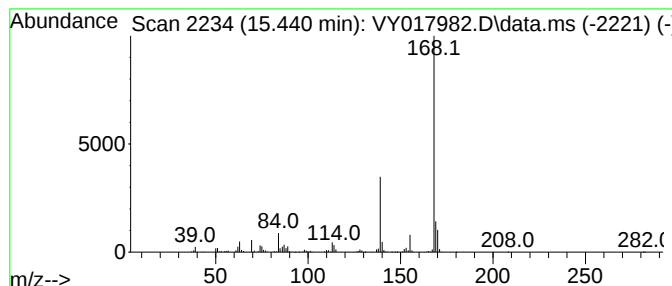
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Quant Title : SW846 8260

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

Peak Number 10 Dibenzofuran Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.440	84.30 ug/l	1578880	1,4-Dichlorobenzene-d4	13.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzofuran	168	C12H8O	000132-64-9	93
2			Tioxolone	168	C7H4O3S	004991-65-5	59
3			2-Amino-6-fluorobenzothiazole	168	C7H5FN2S	000348-40-3	58
4			Naphtho[2,1-d]imidazole	168	C11H8N2	000233-53-4	53
5			9H-Pyrido[3,4-b]indole	168	C11H8N2	000244-63-3	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041824\
Data File : VY017982.D
Acq On : 18 Apr 2024 17:38
Operator : SY/MD
Sample : P2178-01
Misc : 4.17g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TOPSOIL-LKWD-DENALI-4

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041624S.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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Naphthalene, 2,...	12.416	161.9	ug/l	4685130	3	11.495	1446770	50.0
Naphthalene, 1,...	12.971	347.0	ug/l	6498890	4	13.434	936466	50.0
Naphthalene, 1,...	13.111	180.2	ug/l	3375770	4	13.434	936466	50.0
Naphthalene, 2,...	13.580	32.7	ug/l	611848	4	13.434	936466	50.0
Naphthalene, 2,...	13.666	73.3	ug/l	1372150	4	13.434	936466	50.0
Naphthalene, 1,...	14.038	24.6	ug/l	461046	4	13.434	936466	50.0
1,1'-Biphenyl, ...	14.306	21.9	ug/l	409989	4	13.434	936466	50.0
3-Fluoro-para-a...	14.739	269.1	ug/l	5041040	4	13.434	936466	50.0
Dibenzofuran	15.440	84.3	ug/l	1578880	4	13.434	936466	50.0