

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
 Data File : BP019214.D
 Acq On : 02 Jan 2024 18:37
 Operator : MA/JU
 Sample : 05919-15
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCF47

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % 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:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP019214.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.675	97	100	107	rBV	42587	72242	0.14%	0.022%
2	3.763	111	115	123	rVB	50214	72753	0.14%	0.022%
3	4.899	303	308	318	rVB3	42897	88313	0.17%	0.026%
4	5.005	321	326	337	rVV2	37277	70371	0.14%	0.021%
5	6.981	657	662	679	rVB	502803	857448	1.65%	0.256%
6	7.140	683	689	696	rBV	400462	625629	1.21%	0.187%
7	7.328	715	721	731	rBV	514375	824484	1.59%	0.246%
8	7.799	792	801	808	rBV	796777	1262610	2.43%	0.377%
9	7.869	808	813	827	rBV	112182	234863	0.45%	0.070%
10	8.522	917	924	934	rBV	601729	1012374	1.95%	0.302%
11	8.963	992	999	1007	rBV	472447	790006	1.52%	0.236%
12	9.681	1114	1121	1131	rBV	416152	688361	1.33%	0.206%
13	10.010	1170	1177	1186	rBV	225341	411263	0.79%	0.123%
14	10.222	1207	1213	1222	rVB	612873	1037332	2.00%	0.310%
15	10.593	1266	1276	1283	rVB	1020893	1774172	3.42%	0.530%
16	10.710	1292	1296	1300	rBV5	40650	59776	0.12%	0.018%
17	10.751	1300	1303	1309	rVB2	46360	66341	0.13%	0.020%
18	11.075	1353	1358	1367	rBV3	104622	171607	0.33%	0.051%
19	11.269	1384	1391	1397	rBV6	83984	212712	0.41%	0.064%
20	11.416	1411	1416	1420	rBV6	65758	139242	0.27%	0.042%
21	11.481	1423	1427	1431	rVV5	111193	182995	0.35%	0.055%
22	11.622	1445	1451	1458	rBV2	353698	834723	1.61%	0.249%
23	11.775	1474	1477	1485	rVB3	103933	186451	0.36%	0.056%
24	11.869	1486	1493	1498	rBV2	220273	409078	0.79%	0.122%
25	12.016	1515	1518	1522	rBV2	84057	134882	0.26%	0.040%
26	12.116	1531	1535	1538	rBV2	176828	246972	0.48%	0.074%
27	12.157	1539	1542	1546	rVV4	138172	194419	0.37%	0.058%
28	12.328	1566	1571	1577	rBV8	134687	407366	0.78%	0.122%
29	12.387	1577	1581	1584	rBV3	205187	303239	0.58%	0.091%
30	12.681	1628	1631	1635	rBV3	198098	259107	0.50%	0.077%
31	12.787	1646	1649	1653	rBV4	145052	237961	0.46%	0.071%
32	12.975	1676	1681	1686	rBV	1775237	2525885	4.87%	0.754%
33	13.022	1686	1689	1695	rVB4	241289	395428	0.76%	0.118%
34	13.187	1714	1717	1721	rVB3	493023	608501	1.17%	0.182%
35	13.245	1723	1727	1732	rBV3	589498	1085204	2.09%	0.324%
36	13.357	1743	1746	1751	rVB3	301319	347281	0.67%	0.104%

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37	13.663	1794	1798	1801	rBV2	628908	969938	1.87%	0.290%
38	13.857	1828	1831	1836	rVB2	1181882	1502082	2.89%	0.449%
39	13.928	1839	1843	1847	rVB	3691335	4949700	9.54%	1.478%
40	14.134	1875	1878	1883	rBV	1701798	2288406	4.41%	0.683%
41	14.357	1909	1916	1920	rBV	11060348	19394225	37.36%	5.792%
42	14.445	1928	1931	1938	rVB2	1987296	3216244	6.20%	0.961%
43	14.857	1998	2001	2004	rVB	1131694	1249080	2.41%	0.373%
44	14.969	2016	2020	2023	rBV2	2629569	3986348	7.68%	1.190%
45	15.045	2030	2033	2035	rBV	1418241	1980088	3.81%	0.591%
46	15.086	2035	2040	2046	rVB	10353064	16652730	32.08%	4.973%
47	15.292	2070	2075	2079	rBV	4292776	6979105	13.45%	2.084%
48	15.592	2116	2126	2135	rVB	18165602	50947906	98.15%	15.215%
49	15.722	2143	2148	2153	rVB	5927770	9766744	18.82%	2.917%
50	16.210	2226	2231	2237	rVB	9393937	16989685	32.73%	5.074%
51	16.910	2339	2350	2351	rBV	15844196	48906973	94.22%	14.606%
52	23.657	3490	3497	3505	rBV	13835873	30349405	58.47%	9.064%
53	23.786	3514	3519	3524	rVB	3148794	5313972	10.24%	1.587%
54	24.004	3549	3556	3565	rBV2	2892371	11012617	21.22%	3.289%
55	25.198	3745	3759	3763	rBV3	13037375	51906029	100.00%	15.501%
56	25.462	3791	3804	3809	rBV2	6827276	27657628	53.28%	8.260%

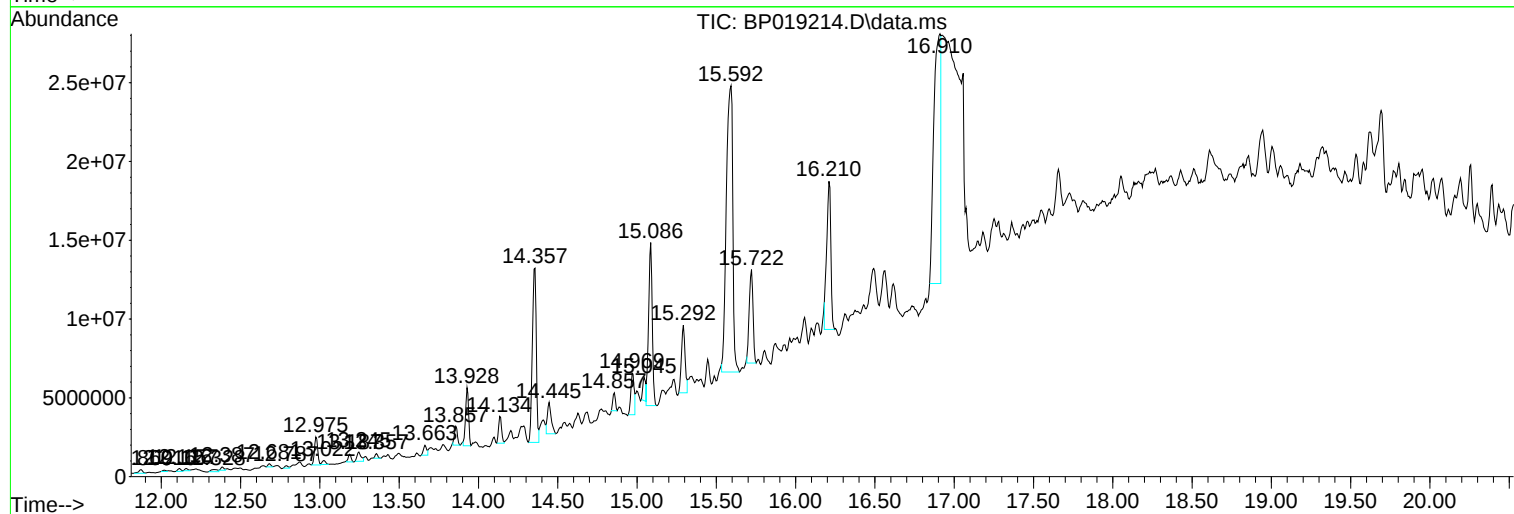
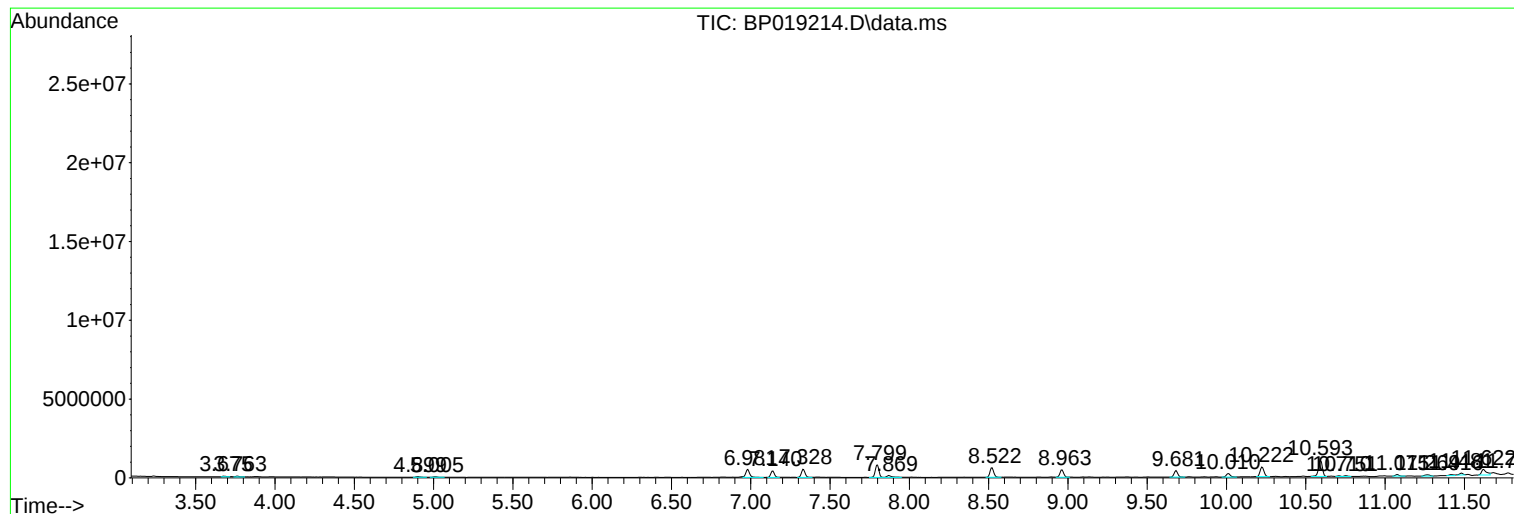
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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



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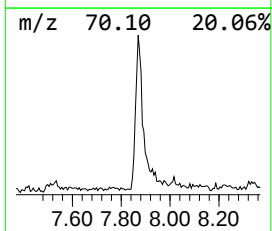
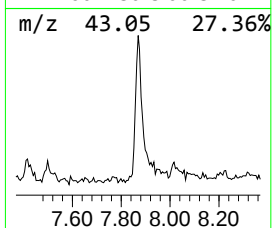
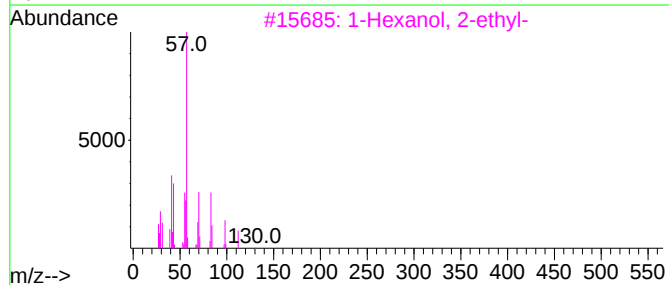
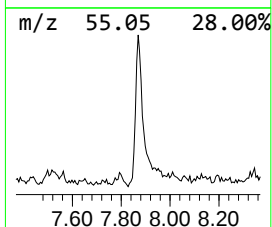
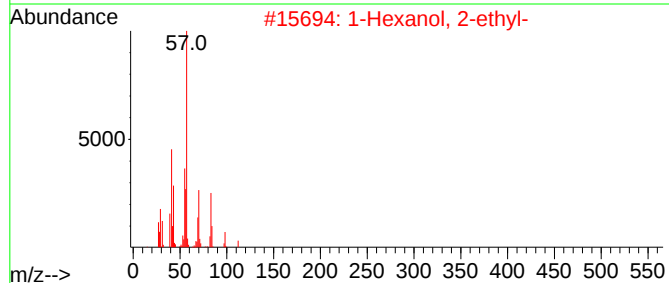
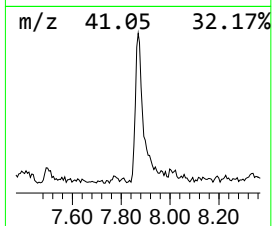
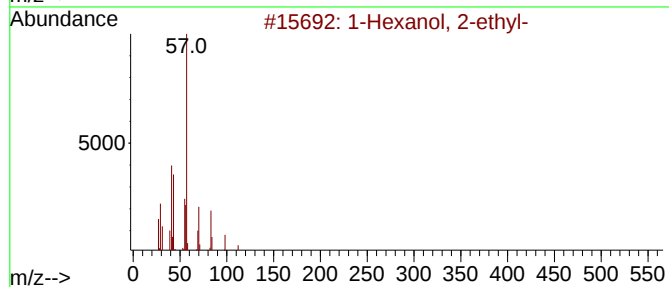
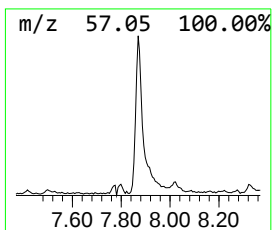
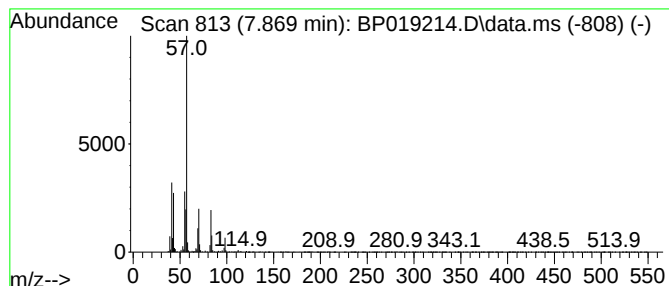
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 1-Hexanol, 2-ethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.869	3.72 ng/ul	234863	1,4-Dichlorobenzene-d4	7.799

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	83
2	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	78
3	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	72
4	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	72
5	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	64



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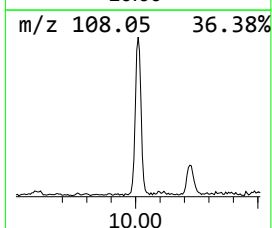
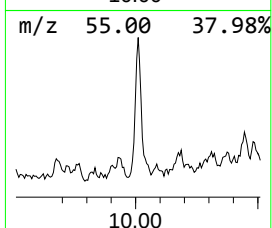
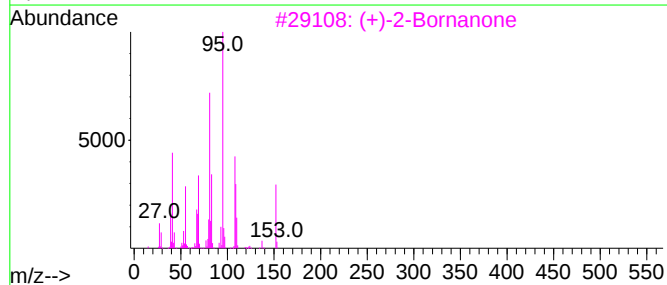
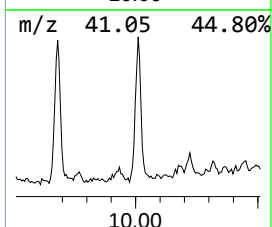
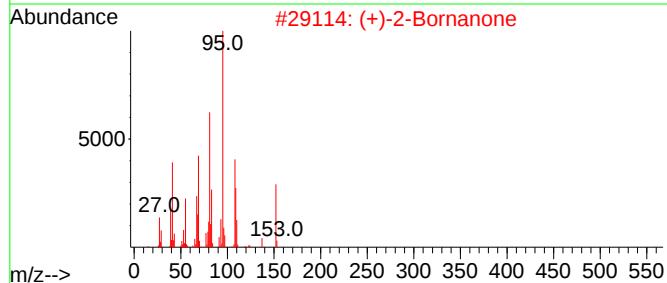
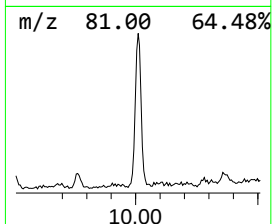
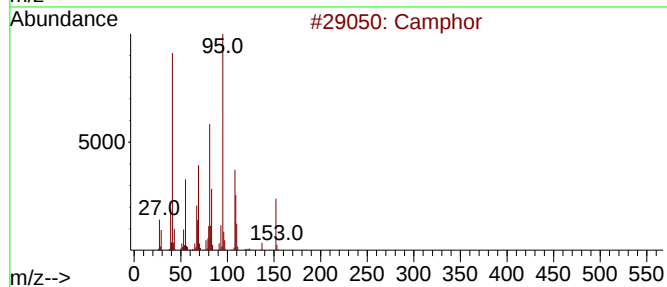
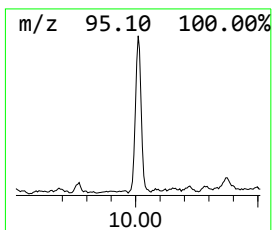
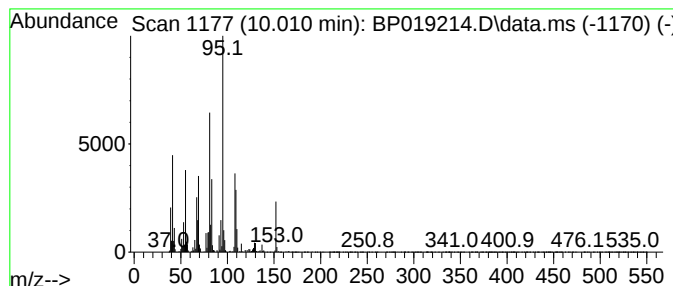
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Camphor Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.010	4.64 ng/ul	411263	Naphthalene-d8	10.592

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Camphor	152	C10H16O	000076-22-2	97
2			(+)-2-Bornanone	152	C10H16O	000464-49-3	97
3			(+)-2-Bornanone	152	C10H16O	000464-49-3	97
4			(+)-2-Bornanone	152	C10H16O	000464-49-3	96
5			(+)-2-Bornanone	152	C10H16O	000464-49-3	96



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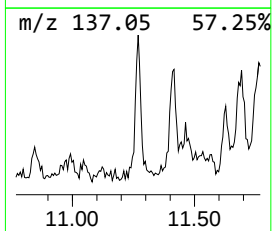
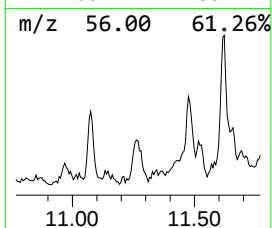
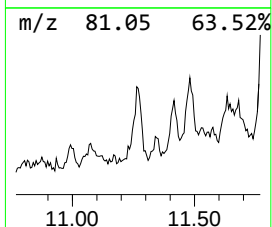
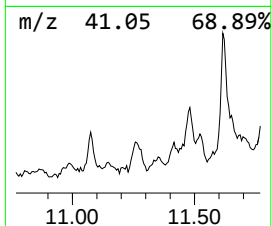
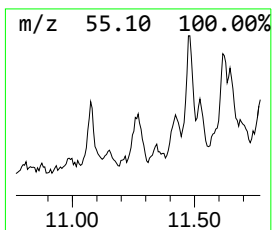
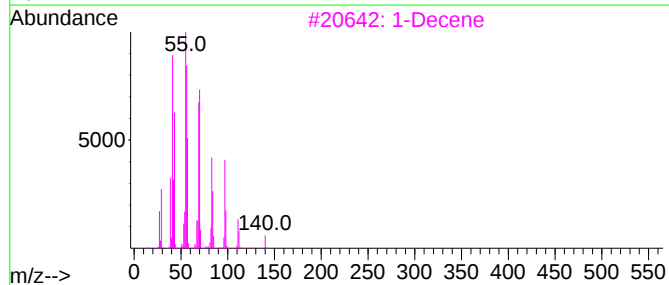
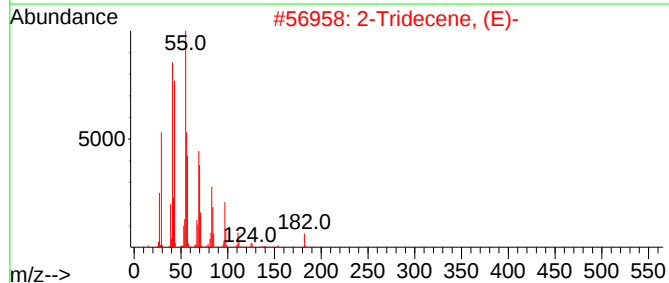
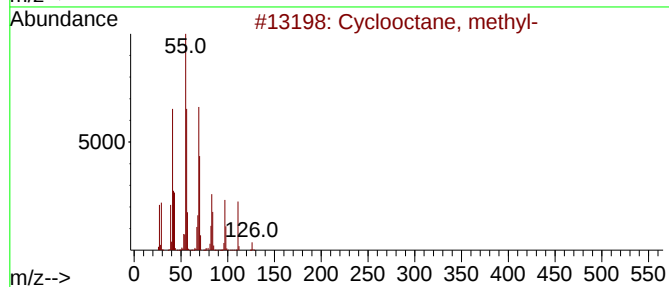
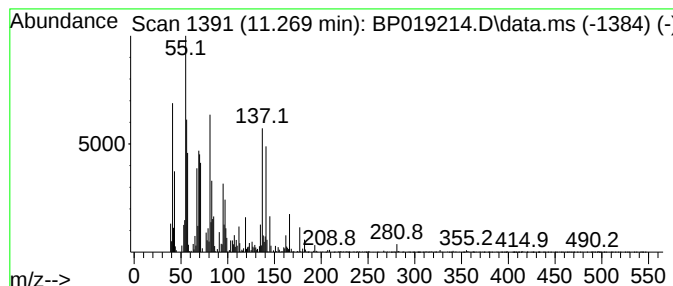
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Cyclic11.269 Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.269	2.40 ng/ul	212712	Naphthalene-d8	10.592

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclooctane, methyl-	126	C9H18	001502-38-1	38
2			2-Tridecene, (E)-	182	C13H26	041446-58-6	38
3			1-Decene	140	C10H20	000872-05-9	35
4			5-Tetradecene, (E)-	196	C14H28	041446-66-6	27
5			cis,trans-2-Ethylbicyclo[4.4.0]d...	166	C12H22	066660-38-6	25



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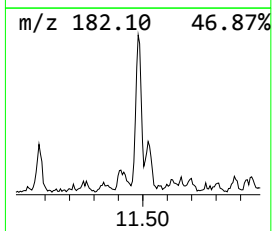
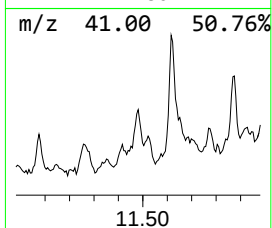
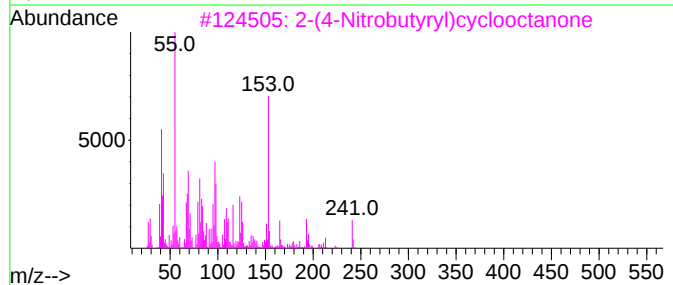
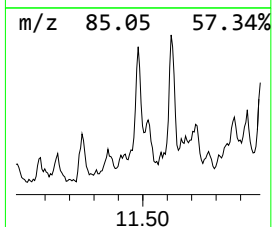
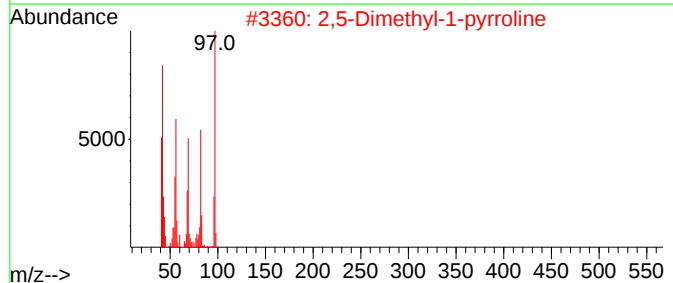
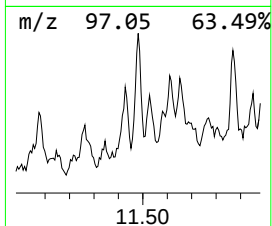
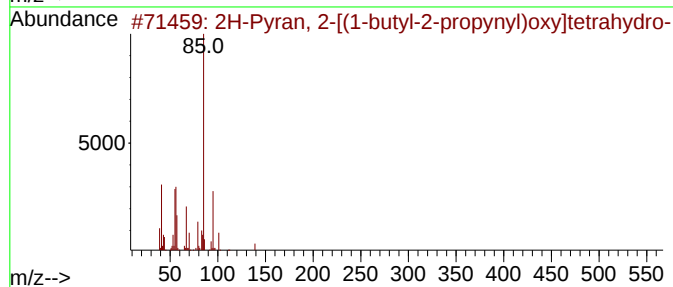
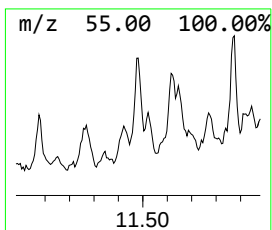
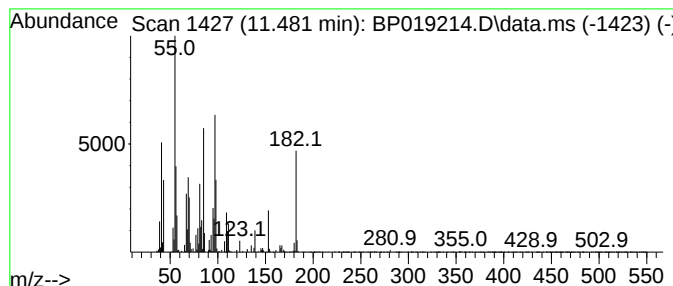
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown-01 Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.481	2.06 ng/ul	182995	Naphthalene-d8	10.592

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Pyran, 2-[(1-butyl-2-propynyl...	196	C12H20O2	000831-83-4	25		
2	2,5-Dimethyl-1-pyrroline	97	C6H11N	000822-64-0	18		
3	2-(4-Nitrobutyryl)cyclooctanone	241	C12H19NO4	079630-93-6	14		
4	Cycloheptane, methyl-	112	C8H16	004126-78-7	14		
5	1H-Pyrazol-4-amine, 3-methyl-	97	C4H7N3	1010338-28-2	14		



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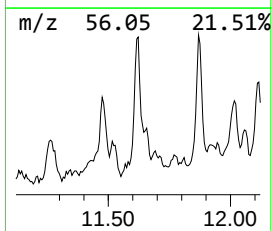
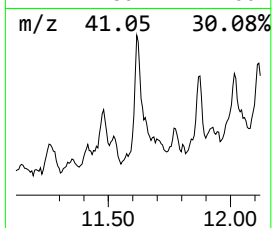
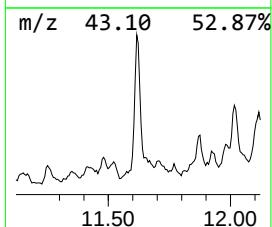
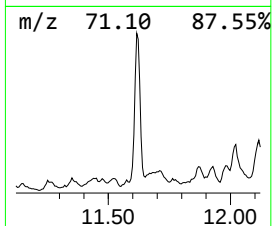
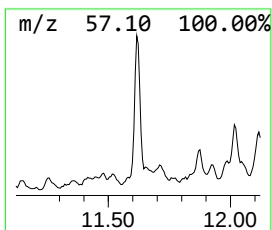
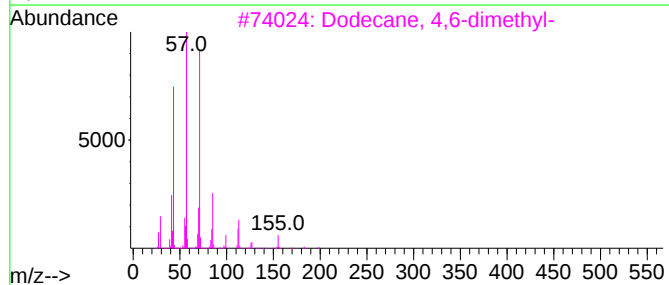
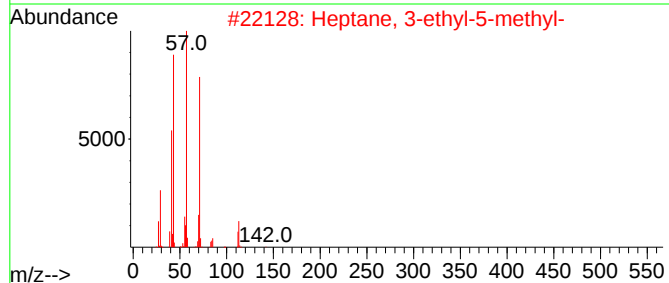
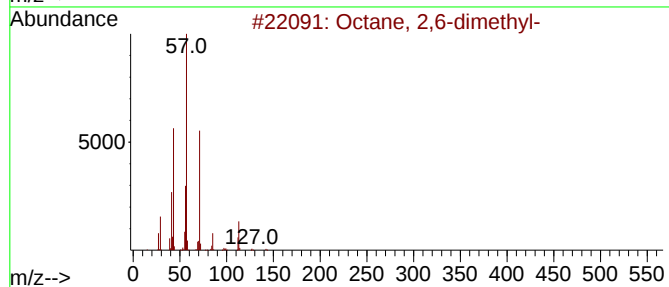
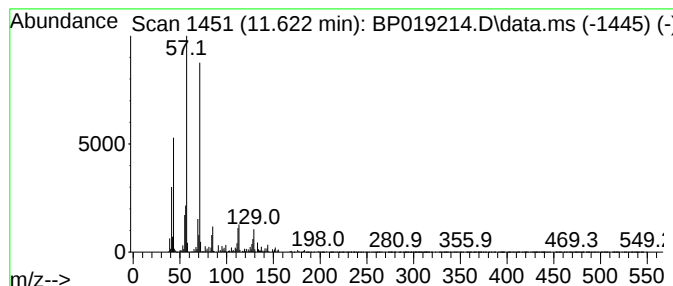
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.622	9.41 ng/ul	834723	Naphthalene-d8	10.592

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	80
2			Heptane, 3-ethyl-5-methyl-	142	C10H22	052896-90-9	64
3			Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	59
4			Sulfurous acid, 2-ethylhexyl pen...	264	C13H28O3S	1000309-18-9	53
5			Oxalic acid, 2-ethylhexyl ethyl ...	230	C12H22O4	1000309-38-4	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
Data File : BP019214.D
Acq On : 02 Jan 2024 18:37
Operator : MA/JU
Sample : 05919-15
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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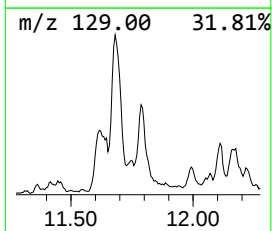
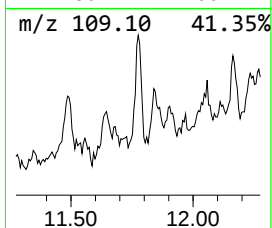
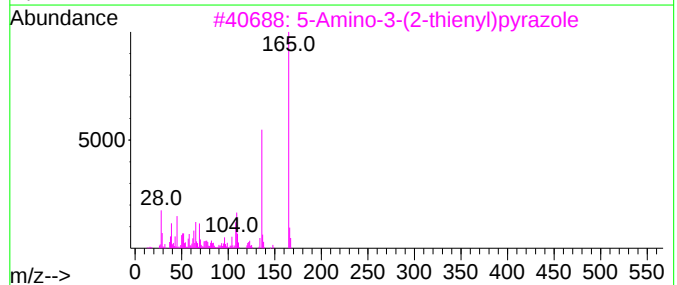
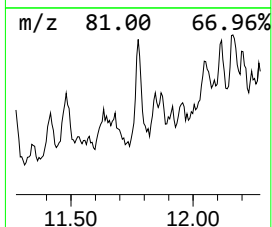
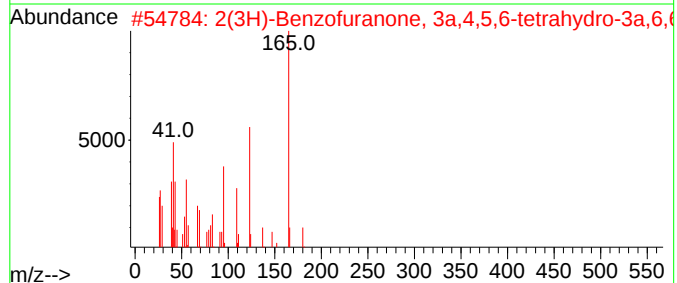
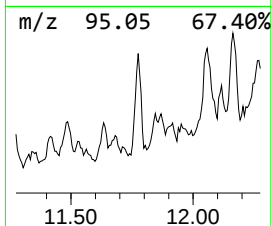
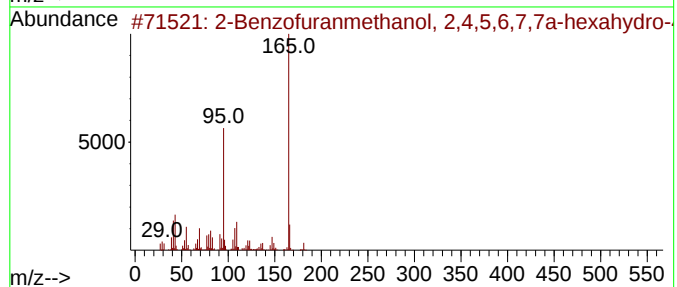
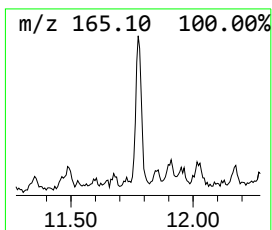
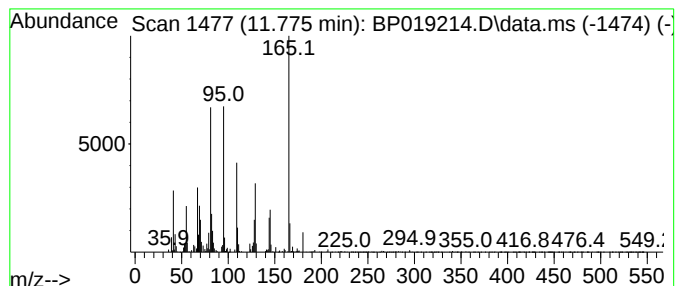
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-02 Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.775	2.10 ng/ul	186451	Naphthalene-d8	10.592

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Benzofuranmethanol, 2,4,5,6,7,...	196	C12H20O2	077384-15-7	38		
2	2(3H)-Benzofuranone, 3a,4,5,6-te...	180	C11H16O2	016778-26-0	38		
3	5-Amino-3-(2-thienyl)pyrazole	165	C7H7N3S	096799-03-0	27		
4	2-Methyl-5-benzothiazolol	165	C8H7NOS	068867-14-1	22		
5	4,6-Dimethyl-7-oxo-4,7-dihydro-t...	165	C6H7N5O	025623-69-2	22		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
Data File : BP019214.D
Acq On : 02 Jan 2024 18:37
Operator : MA/JU
Sample : 05919-15
Misc :
ALS Vial : 15 Sample Multiplier: 1

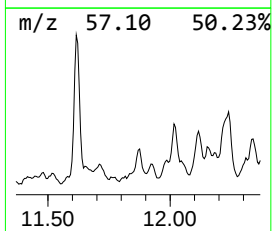
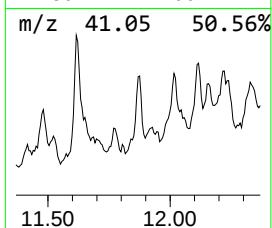
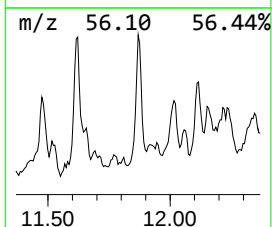
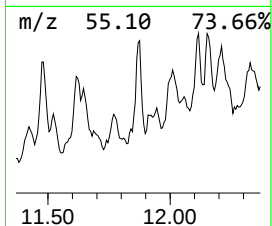
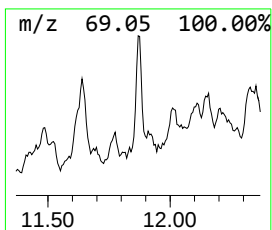
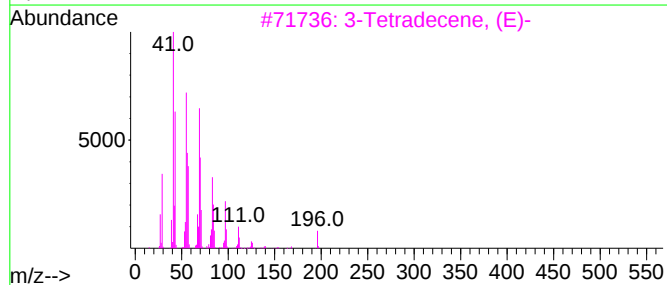
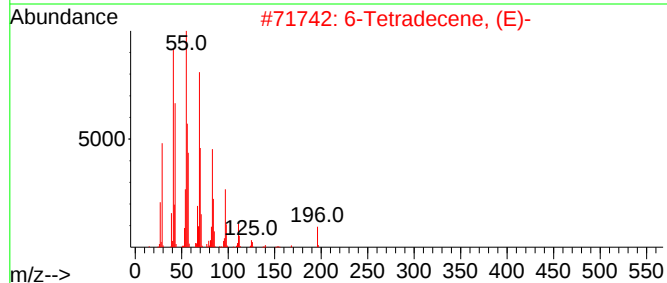
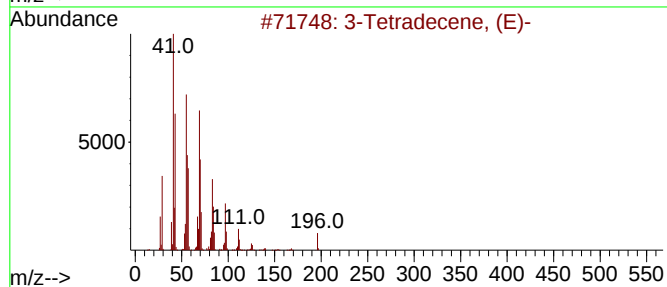
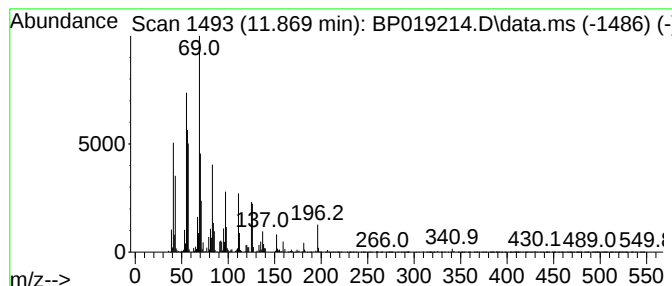
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.869	4.61 ng/ul	409078	Naphthalene-d8	10.592	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Tetradecene, (E)-	196	C14H28	041446-68-8	90
2	6-Tetradecene, (E)-	196	C14H28	041446-64-4	90
3	3-Tetradecene, (E)-	196	C14H28	041446-68-8	90
4	6-Dodecene, (Z)-	168	C12H24	007206-29-3	76
5	6-Tridecene, (E)-	182	C13H26	006434-76-0	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
Data File : BP019214.D
Acq On : 02 Jan 2024 18:37
Operator : MA/JU
Sample : 05919-15
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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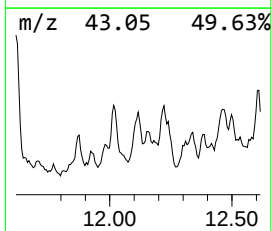
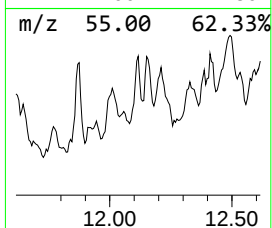
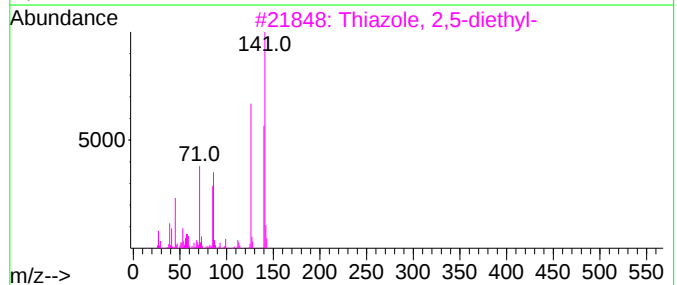
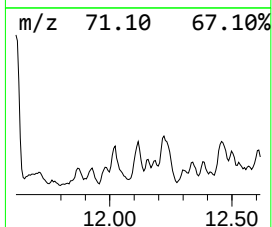
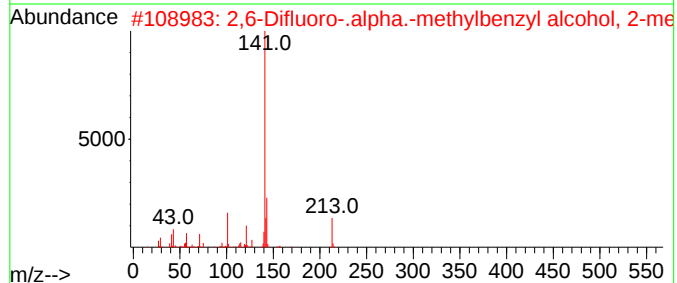
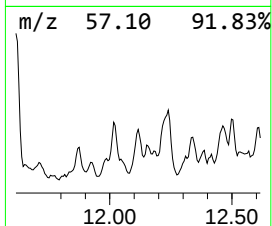
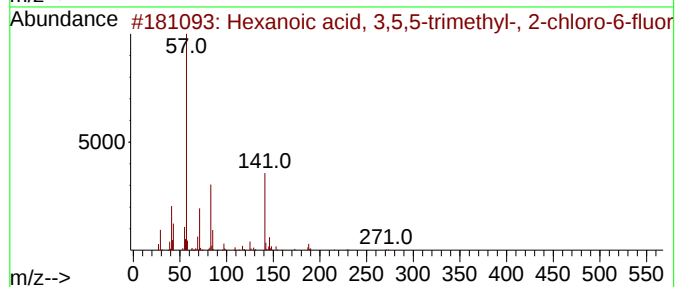
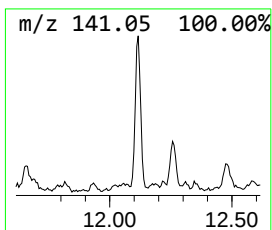
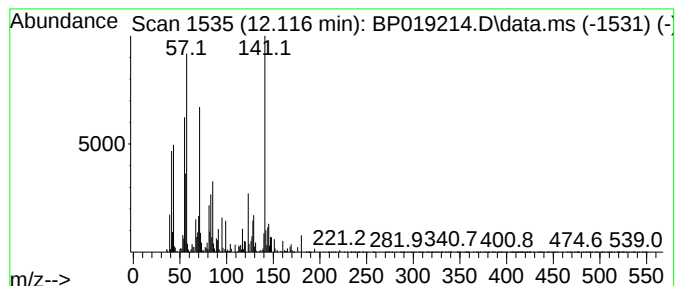
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Hexanoic acid, 3,5,5-trimet... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.116	2.78 ng/ul	246972	Naphthalene-d8	10.592

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 3,5,5-trimethyl-,...	286	C15H20ClF02	1000406-82-4	59
2			2,6-Difluoro-.alpha.-methylbenzy...	228	C13H18F2O	1000395-25-5	47
3			Thiazole, 2,5-diethyl-	141	C7H11NS	015729-76-7	47
4			Nonanoic acid, 3,4-dichloropheny...	302	C15H20Cl2O2	1000358-02-7	47
5			Thiazole, 2,4-diethyl-	141	C7H11NS	032272-49-4	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
Data File : BP019214.D
Acq On : 02 Jan 2024 18:37
Operator : MA/JU
Sample : 05919-15
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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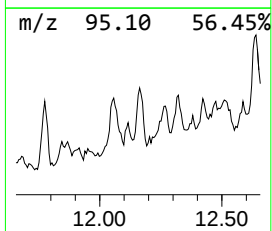
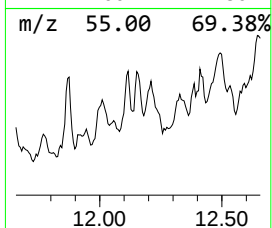
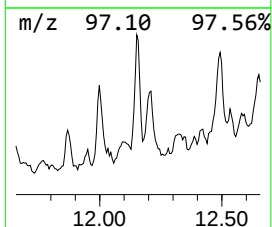
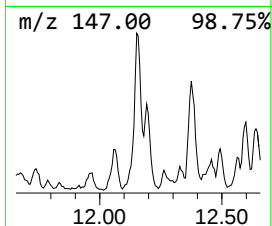
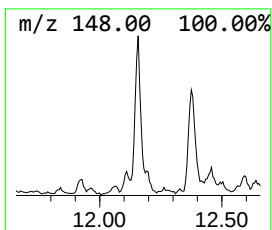
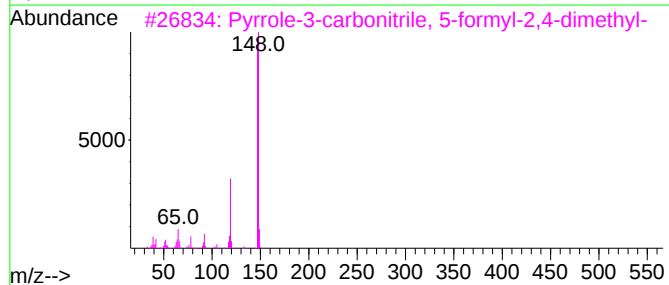
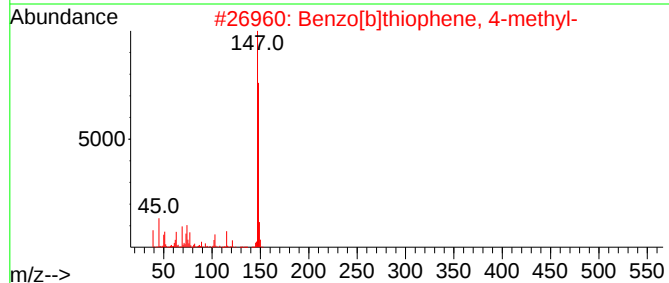
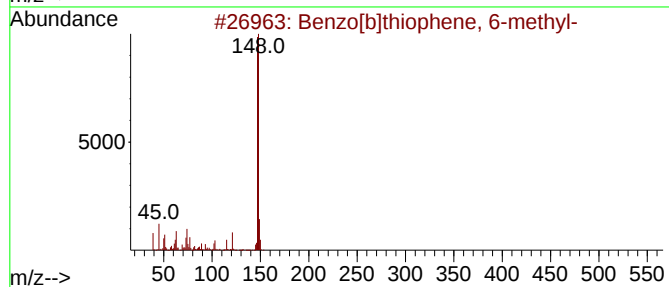
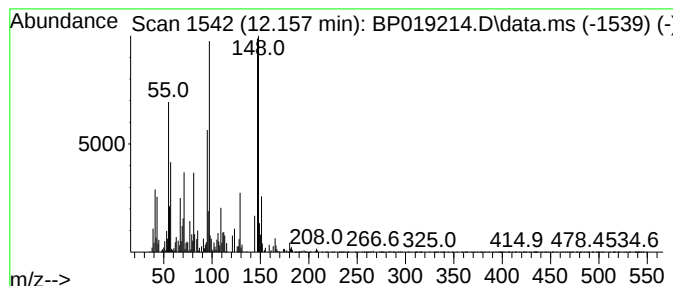
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-03 Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.157	2.19 ng/ul	194419	Naphthalene-d8	10.592

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[b]thiophene, 6-methyl-	148	C9H8S	016587-47-6	38
2			Benzo[b]thiophene, 4-methyl-	148	C9H8S	014315-11-8	35
3			Pyrrole-3-carbonitrile, 5-formyl...	148	C8H8N2O	032487-71-1	35
4			3-Methylbenzothiophene	148	C9H8S	001455-18-1	35
5			2-Methyl-5-hydroxybenzofuran	148	C9H8O2	006769-56-8	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
Data File : BP019214.D
Acq On : 02 Jan 2024 18:37
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Sample : 05919-15
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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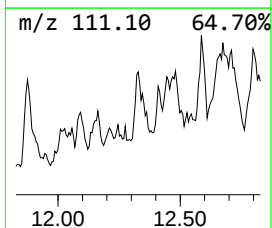
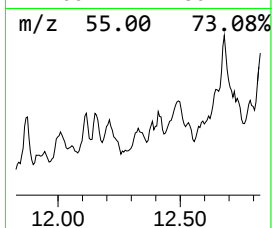
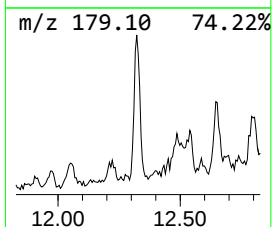
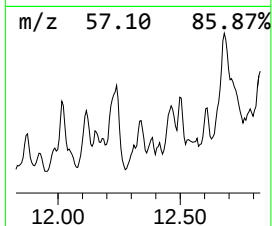
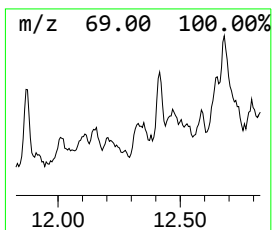
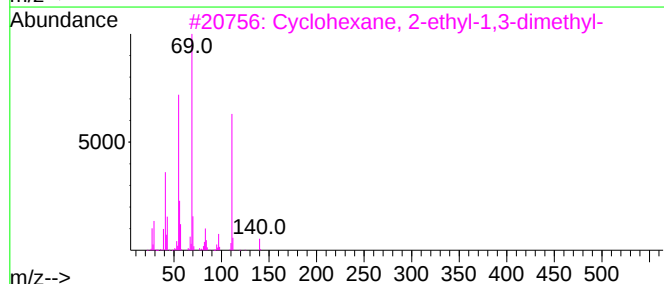
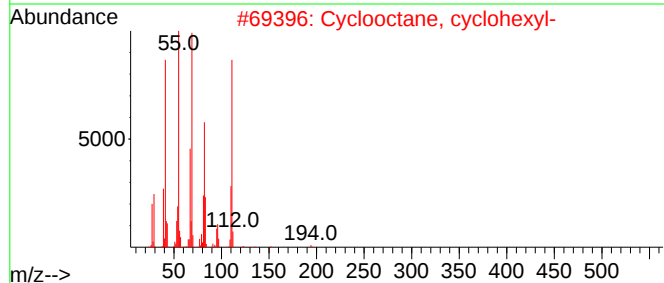
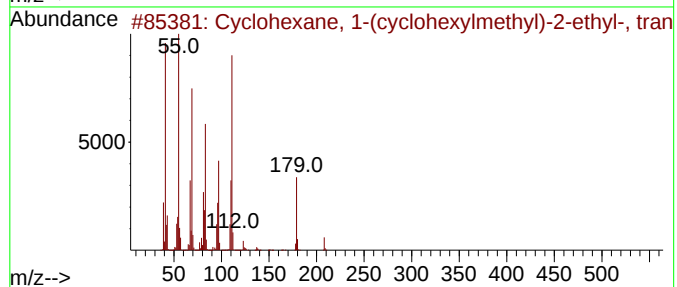
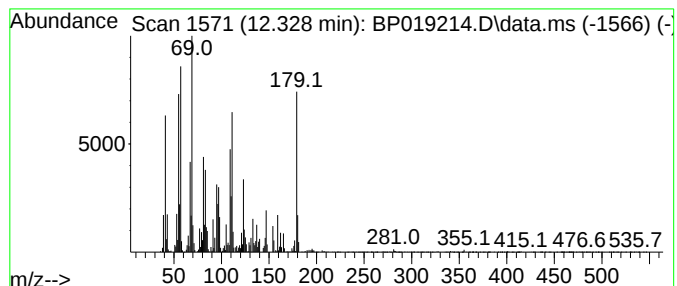
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-04 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.328	4.59 ng/ul	407366	Naphthalene-d8	10.592

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-(cyclohexylmethyl...	208	C15H28	054934-92-8	43
2			Cyclooctane, cyclohexyl-	194	C14H26	092369-78-3	42
3			Cyclohexane, 2-ethyl-1,3-dimethyl-	140	C10H20	007045-67-2	25
4			Cyclooctane, ethyl-	140	C10H20	013152-02-8	25
5			Cyclohexane, 1,2-diethyl-, cis-	140	C10H20	000824-43-1	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
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Acq On : 02 Jan 2024 18:37
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Misc :
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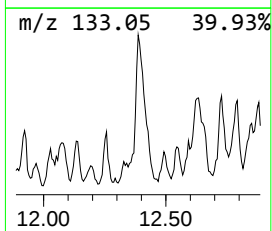
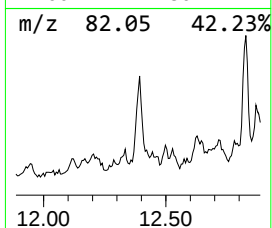
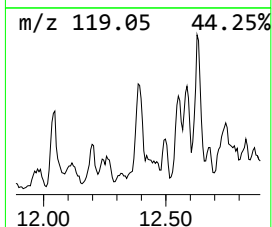
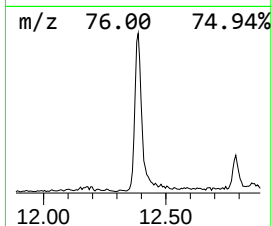
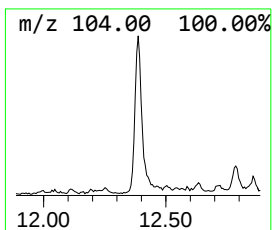
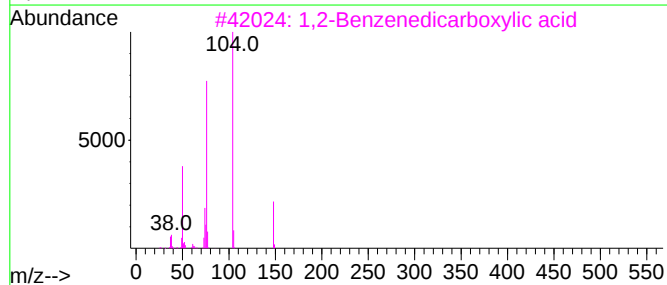
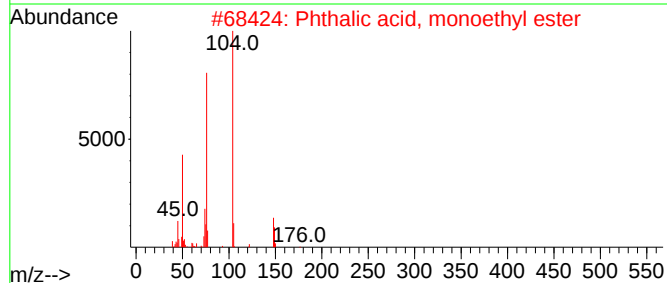
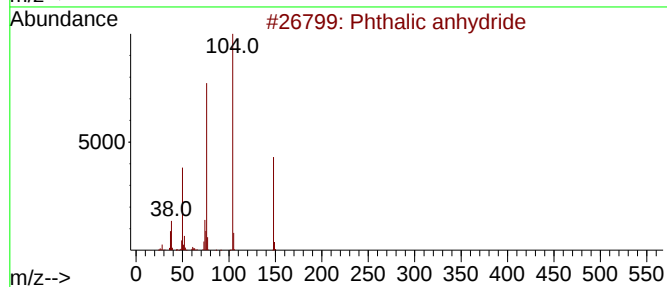
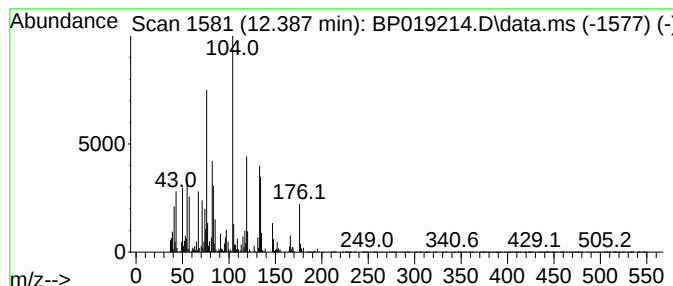
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Phthalic anhydride Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.387	3.42 ng/ul	303239	Naphthalene-d8	10.592

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic anhydride	148	C8H4O3	000085-44-9	64
2			Phthalic acid, monoethyl ester	194	C10H10O4	002306-33-4	50
3			1,2-Benzenedicarboxylic acid	166	C8H6O4	000088-99-3	50
4			Bicyclo[4.2.0]octa-1,3,5-triene...	132	C8H4O2	006383-11-5	50
5			Indan-1,2,3-trione	160	C9H4O3	000938-24-9	50



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Data File : BP019214.D
Acq On : 02 Jan 2024 18:37
Operator : MA/JU
Sample : 05919-15
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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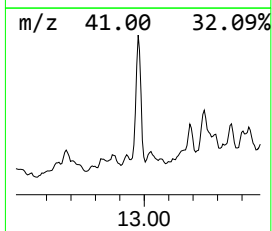
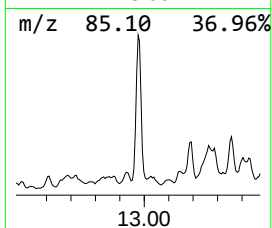
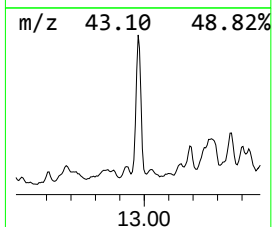
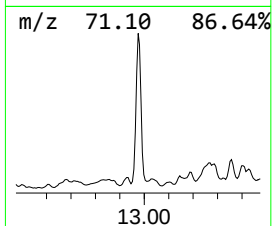
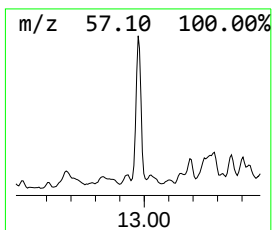
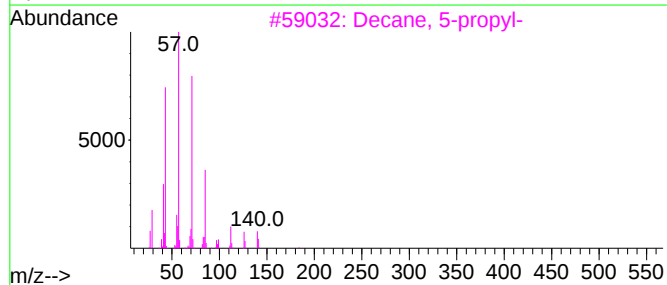
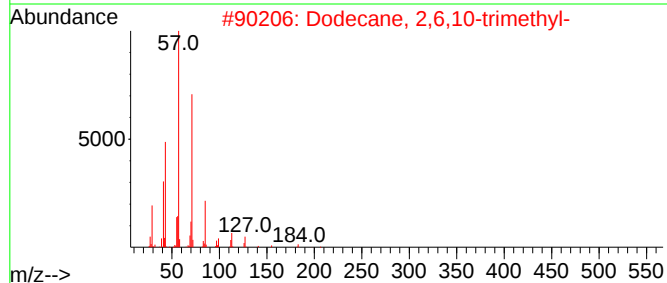
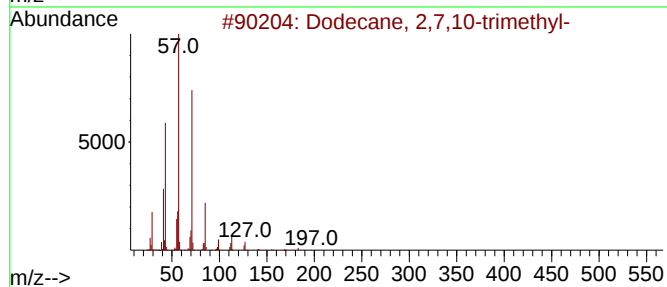
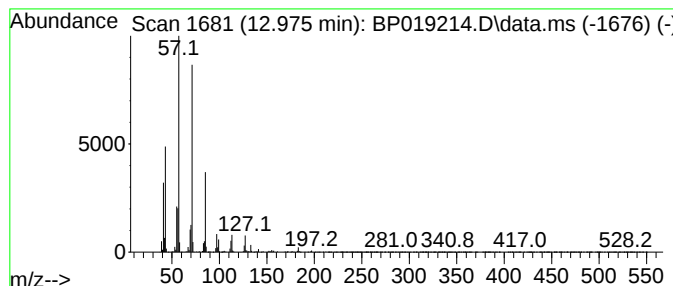
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.975	15.71 ng/ul	2525890	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	91
2			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	90
3			Decane, 5-propyl-	184	C13H28	017312-62-8	80
4			Pentadecane, 7-methyl-	226	C16H34	006165-40-8	78
5			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
Data File : BP019214.D
Acq On : 02 Jan 2024 18:37
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Sample : 05919-15
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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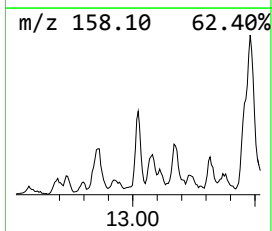
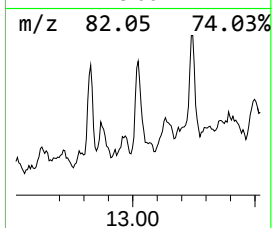
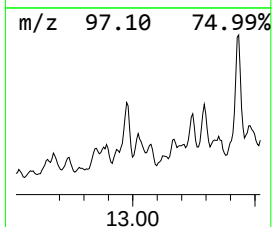
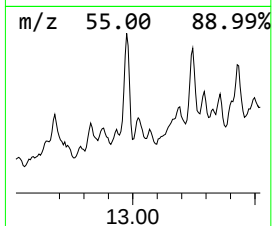
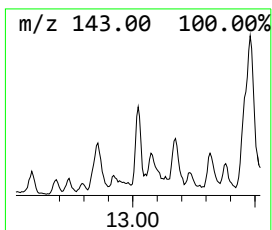
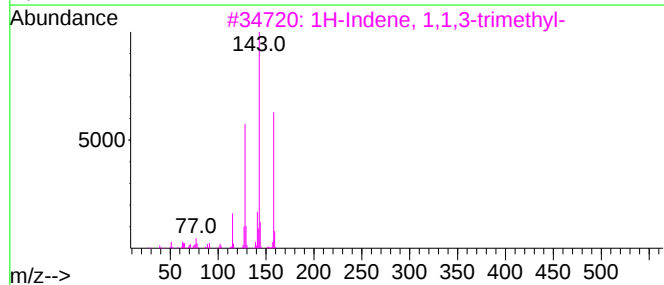
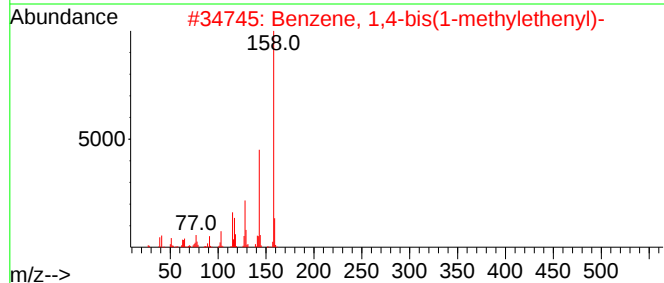
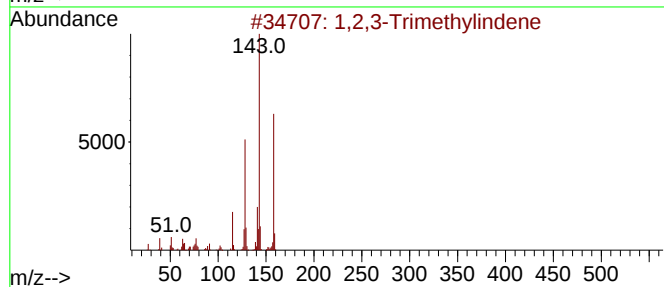
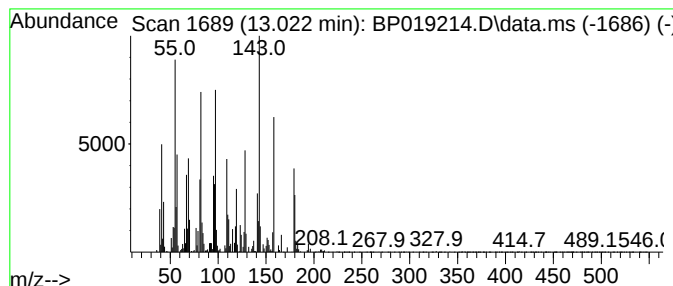
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 unknown-05 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.022	2.46 ng/ul	395428	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2,3-Trimethylindene	158	C12H14	004773-83-5	38
2			Benzene, 1,4-bis(1-methylethenyl)-	158	C12H14	001605-18-1	38
3			1H-Indene, 1,1,3-trimethyl-	158	C12H14	002177-45-9	30
4			(1-Methylpenta-2,4-dienyl)benzene	158	C12H14	1000210-01-1	30
5			1,2,3-Trimethylindene	158	C12H14	004773-83-5	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
Data File : BP019214.D
Acq On : 02 Jan 2024 18:37
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Sample : 05919-15
Misc :
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ClientSampleId :
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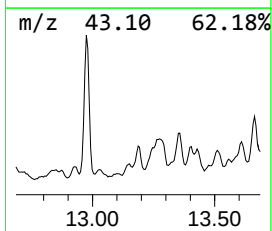
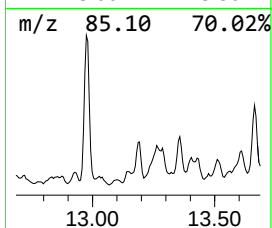
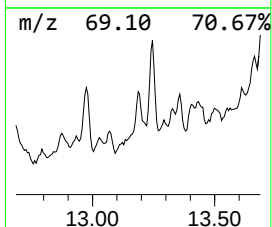
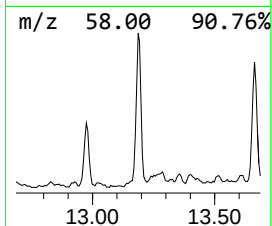
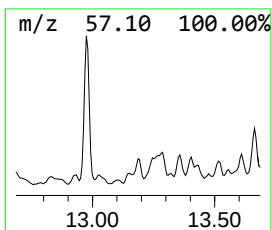
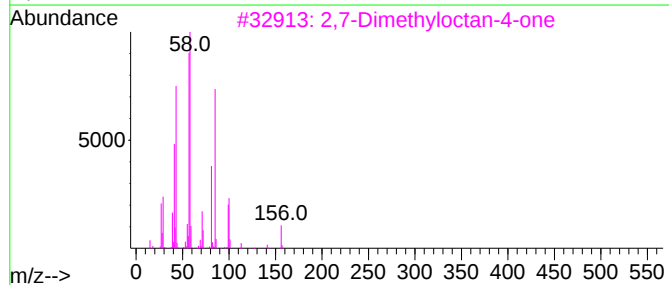
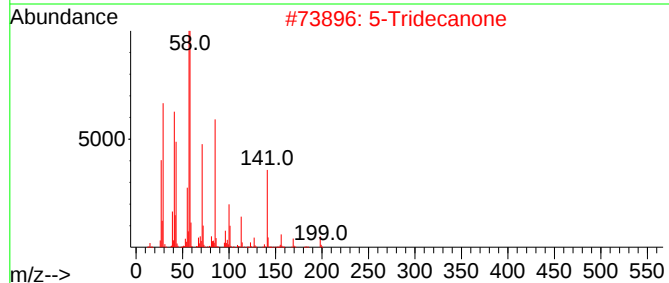
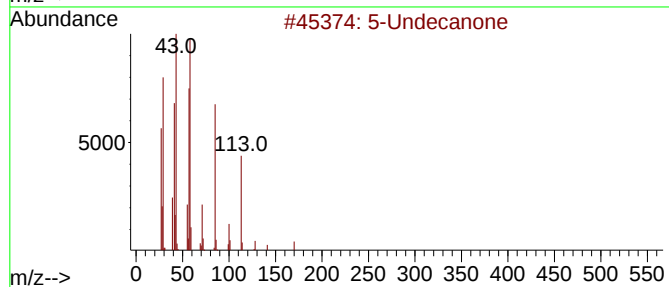
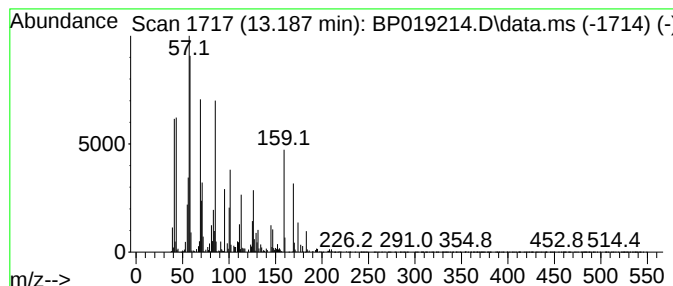
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 unknown-06 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.187	3.78 ng/ul	608501	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Undecanone	170	C11H22O	033083-83-9	38
2			5-Tridecanone	198	C13H26O	030692-16-1	35
3			2,7-Dimethyloctan-4-one	156	C10H20O	059387-92-7	35
4			2-Methyl-4-decanone	170	C11H22O	006628-25-7	35
5			Triallylphosphine	154	C9H15P	016523-89-0	20



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
Data File : BP019214.D
Acq On : 02 Jan 2024 18:37
Operator : MA/JU
Sample : 05919-15
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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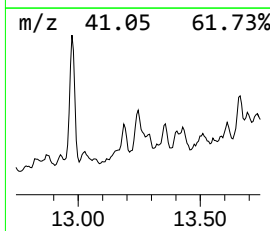
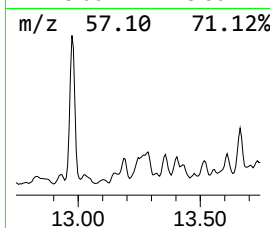
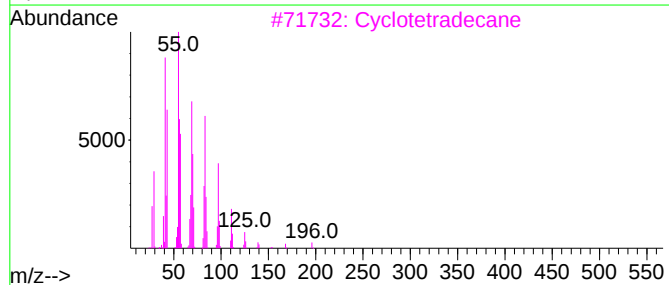
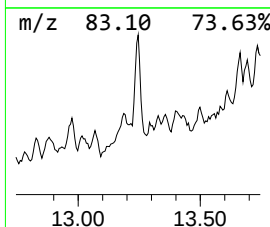
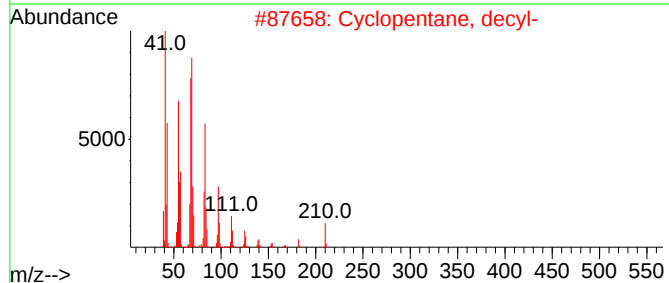
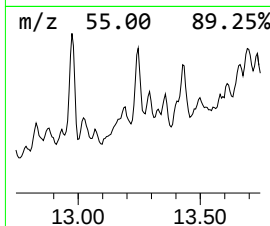
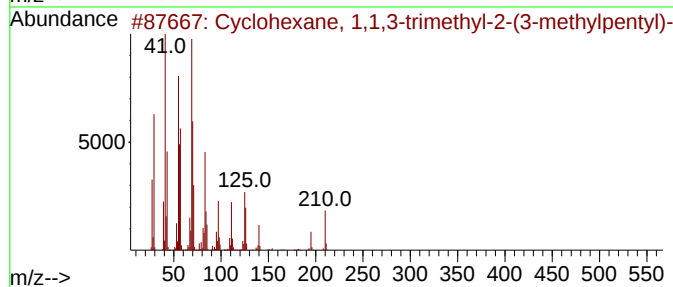
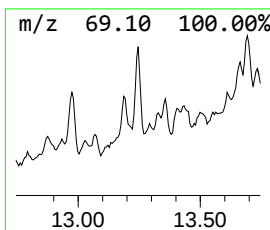
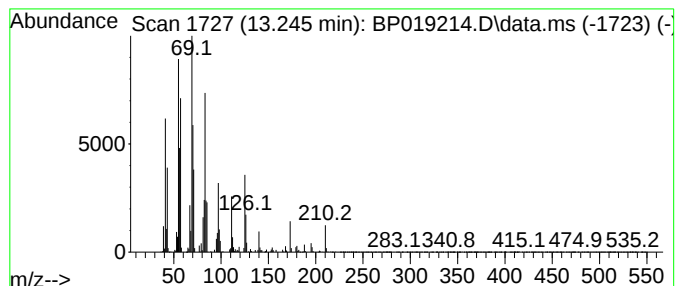
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 (DEL) Alkane: Cyclic13.245 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.245	6.75 ng/ul	1085200	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,1,3-trimethyl-2-(...	210	C15H30	054965-05-8	91
2			Cyclopentane, decyl-	210	C15H30	001795-21-7	62
3			Cyclotetradecane	196	C14H28	000295-17-0	53
4			Cyclotetradecane	196	C14H28	000295-17-0	50
5			5-Ethyl-1-nonene	154	C11H22	019780-74-6	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
Data File : BP019214.D
Acq On : 02 Jan 2024 18:37
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Sample : 05919-15
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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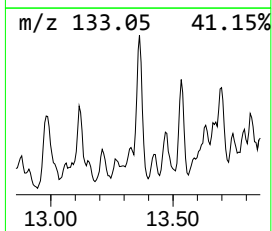
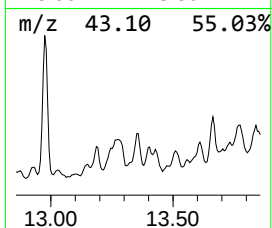
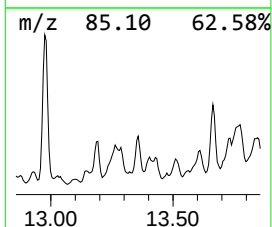
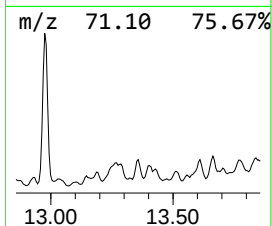
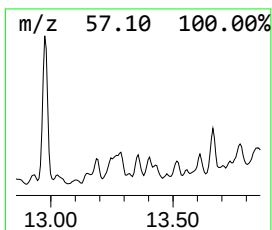
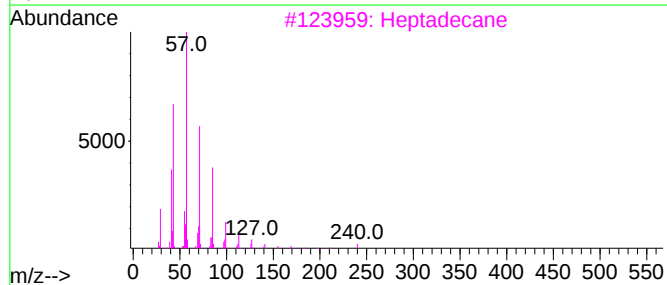
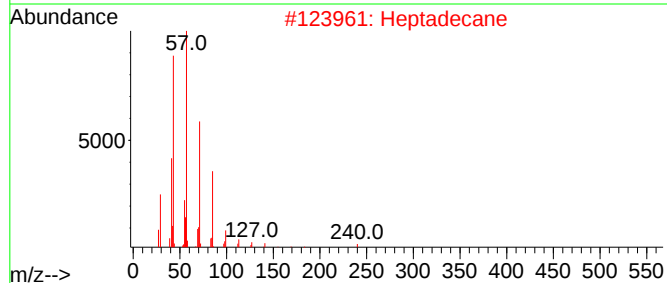
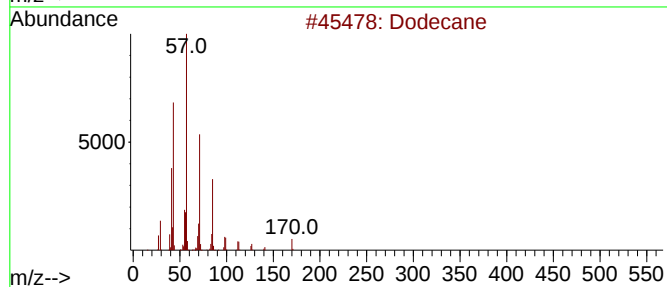
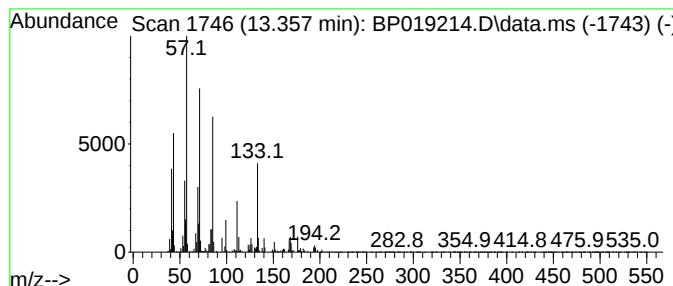
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.357	2.16 ng/ul	347281	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	64
2			Heptadecane	240	C17H36	000629-78-7	62
3			Heptadecane	240	C17H36	000629-78-7	58
4			9-methylheptadecane	254	C18H38	026741-18-4	53
5			Sulfurous acid, dodecyl 2-propyl...	292	C15H32O3S	1000309-12-3	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
Data File : BP019214.D
Acq On : 02 Jan 2024 18:37
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Sample : 05919-15
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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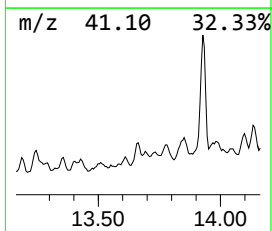
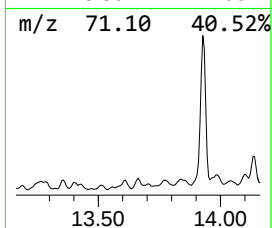
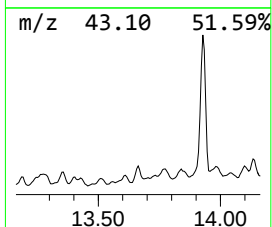
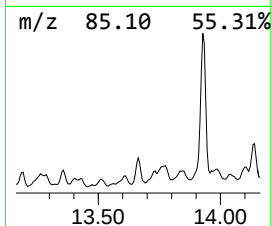
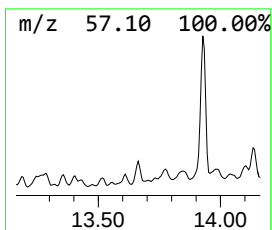
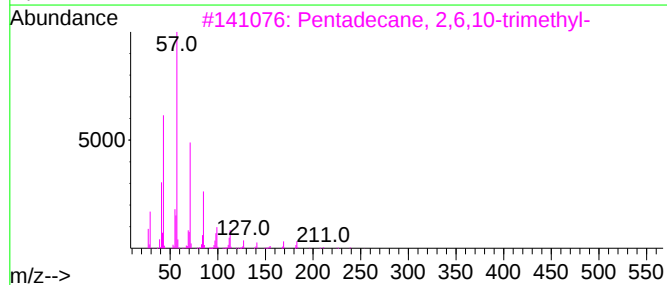
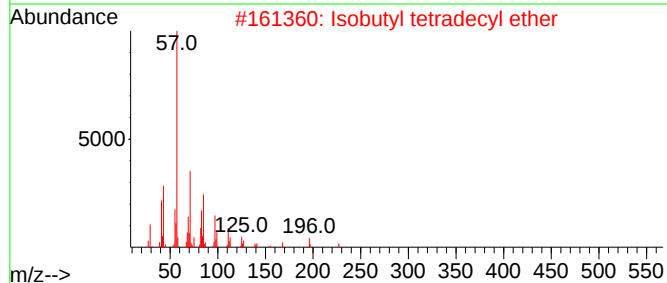
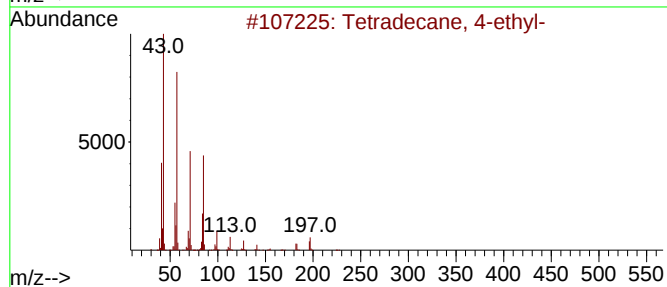
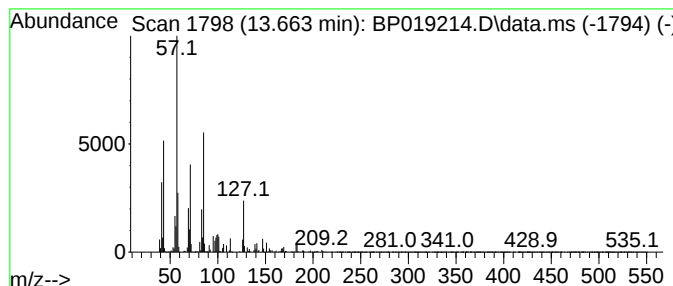
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.663	6.03 ng/ul	969938	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane, 4-ethyl-	226	C16H34	055045-14-2	53
2			Isobutyl tetradecyl ether	270	C18H38O	1000406-32-7	43
3			Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	38
4			Tetradecane	198	C14H30	000629-59-4	38
5			Decane, 3-bromo-	220	C10H21Br	030571-71-2	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
Data File : BP019214.D
Acq On : 02 Jan 2024 18:37
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Sample : 05919-15
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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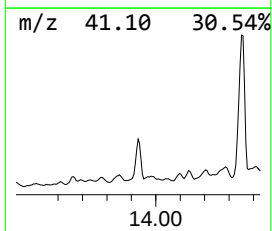
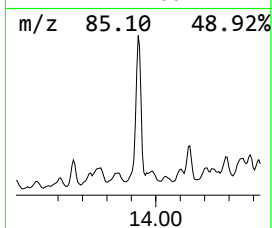
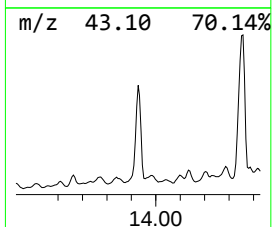
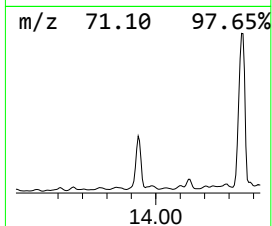
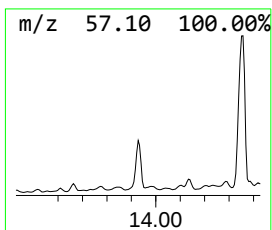
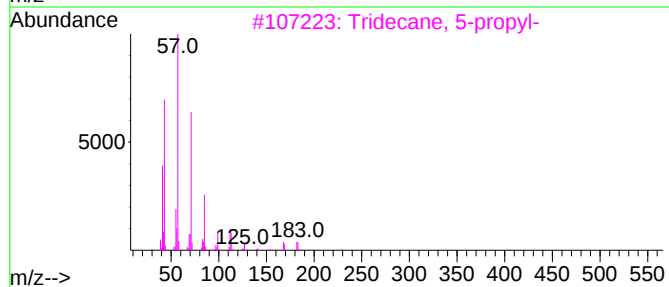
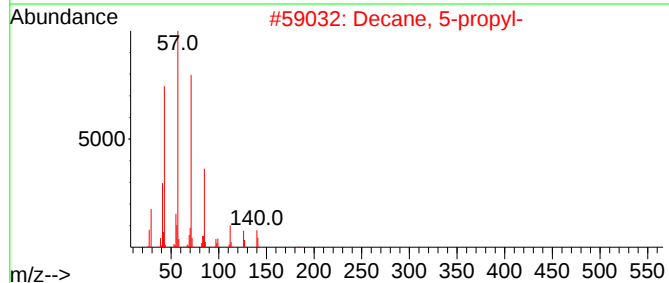
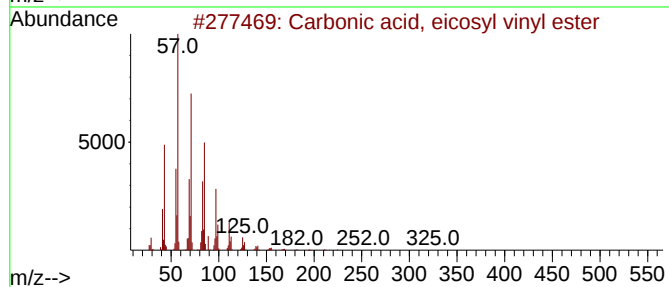
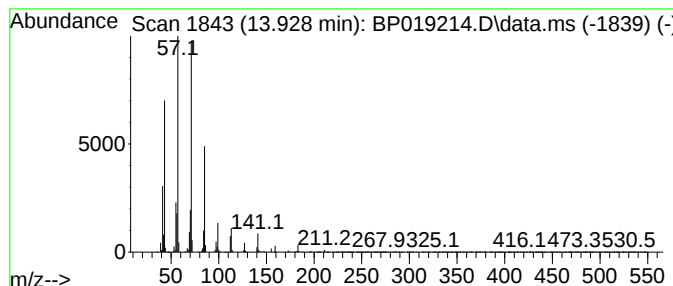
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 Carbonic acid, eicosyl vinyl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.928	30.78 ng/ul	4949700	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	90
2			Decane, 5-propyl-	184	C13H28	017312-62-8	81
3			Tridecane, 5-propyl-	226	C16H34	055045-11-9	76
4			Dodecane, 4-methyl-	184	C13H28	006117-97-1	74
5			Undecane, 4,8-dimethyl-	184	C13H28	017301-33-6	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
Data File : BP019214.D
Acq On : 02 Jan 2024 18:37
Operator : MA/JU
Sample : 05919-15
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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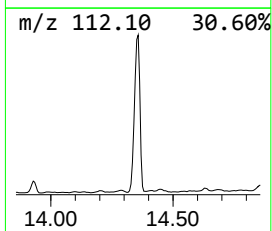
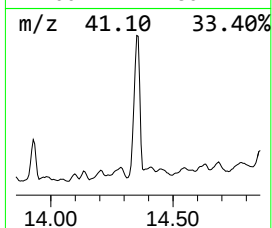
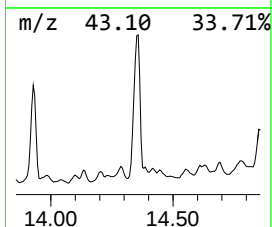
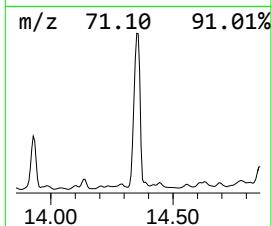
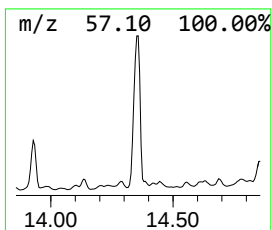
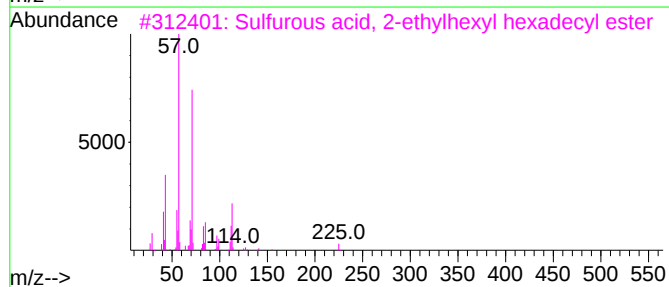
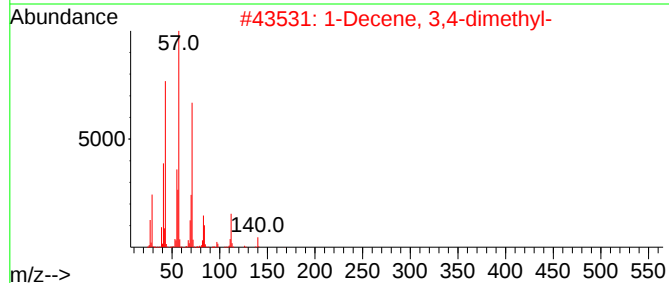
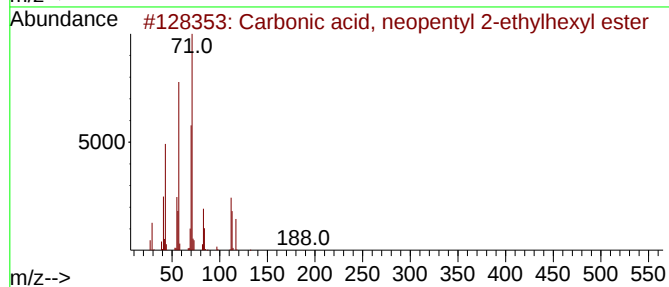
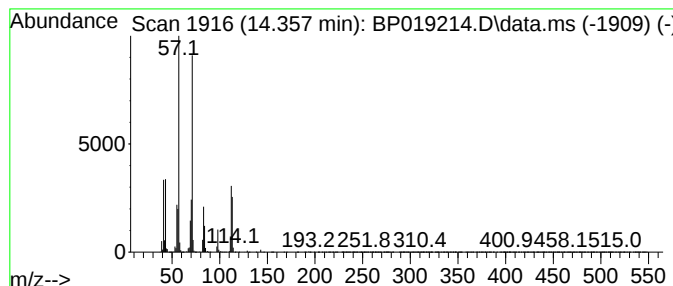
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 Carbonic acid, neopentyl 2-... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.357	120.60 ng/ul	19394200	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, neopentyl 2-ethyl...	244	C14H28O3	1000357-85-7	72
2			1-Decene, 3,4-dimethyl-	168	C12H24	050871-03-9	72
3			Sulfurous acid, 2-ethylhexyl hex...	418	C24H50O3S	1000309-19-9	59
4			Carbonic acid, 2-ethylhexyl nony...	300	C18H36O3	1000383-13-6	59
5			Carbonic acid, octyl vinyl ester	200	C11H20O3	1000383-25-5	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
Data File : BP019214.D
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Operator : MA/JU
Sample : 05919-15
Misc :
ALS Vial : 15 Sample Multiplier: 1

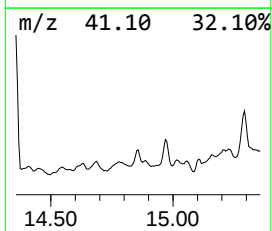
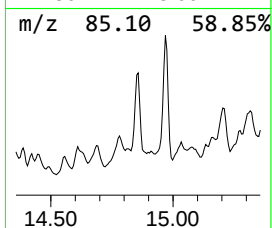
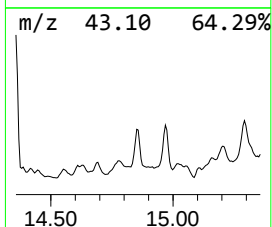
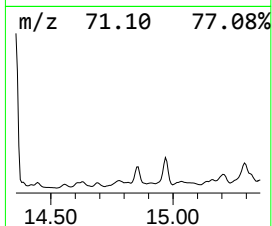
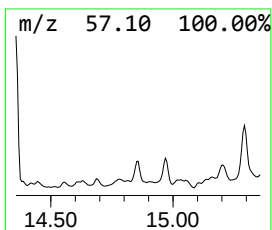
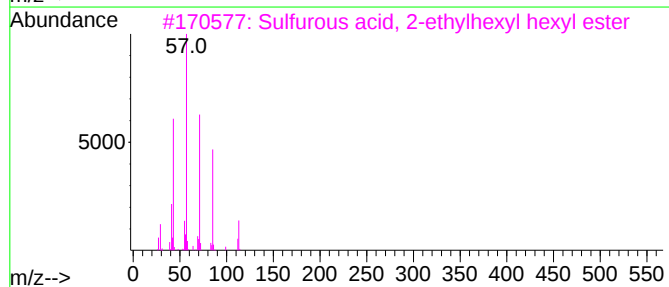
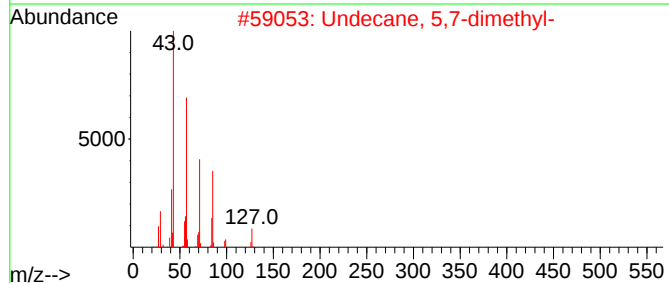
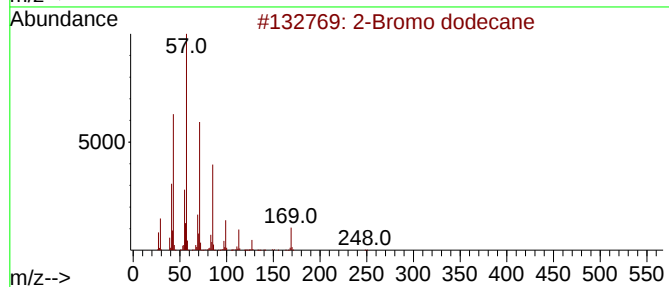
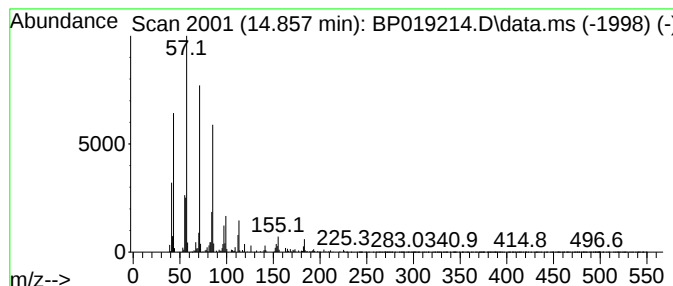
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 2-Bromo dodecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.857	7.77 ng/ul	1249080	Acenaphthene-d10	14.445	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Bromo dodecane	248	C12H25Br	013187-99-0	80
2	Undecane, 5,7-dimethyl-	184	C13H28	017312-83-3	64
3	Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	64
4	Sulfurous acid, 2-ethylhexyl iso...	278	C14H30O3S	1000309-19-0	64
5	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
Data File : BP019214.D
Acq On : 02 Jan 2024 18:37
Operator : MA/JU
Sample : 05919-15
Misc :
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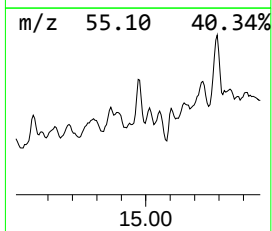
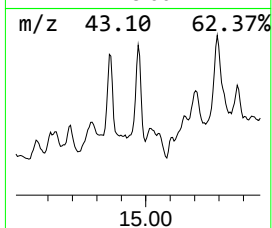
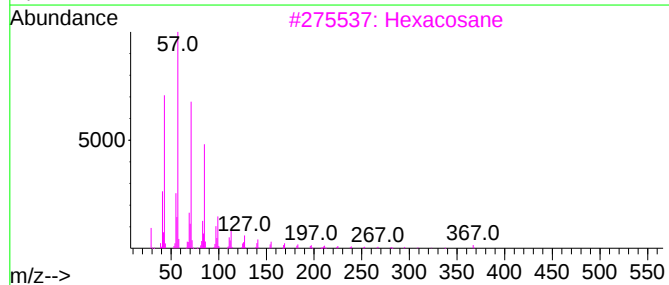
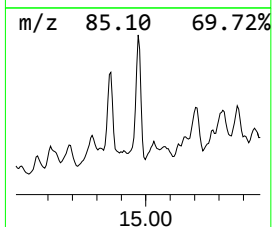
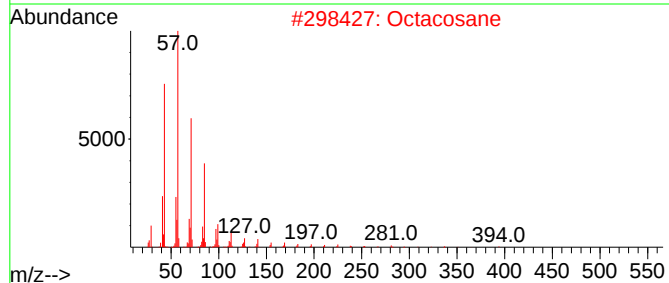
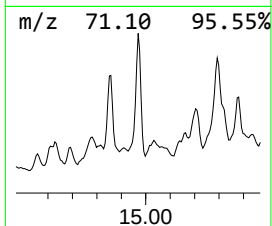
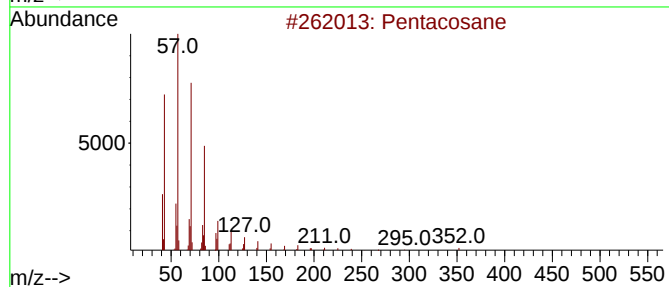
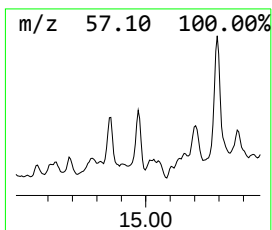
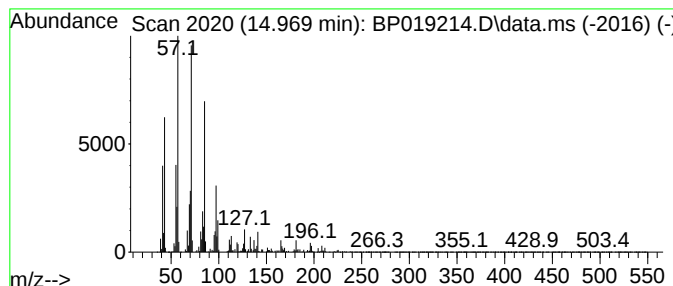
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.969	24.79 ng/ul	3986350	Acenaphthene-d10		14.445	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Pentacosane		352	C25H52	000629-99-2	72
2	Octacosane		394	C28H58	000630-02-4	72
3	Hexacosane		366	C26H54	000630-01-3	72
4	Heptadecane, 8-methyl-		254	C18H38	013287-23-5	72
5	Dodecane, 2-methyl-6-propyl-		226	C16H34	055045-08-4	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
Data File : BP019214.D
Acq On : 02 Jan 2024 18:37
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Sample : 05919-15
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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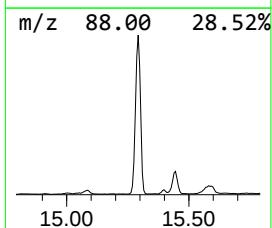
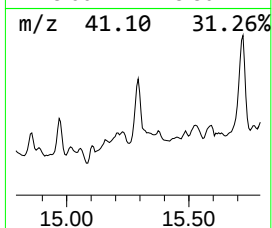
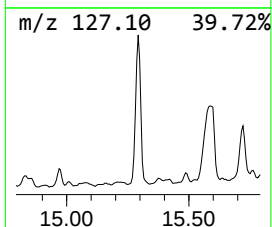
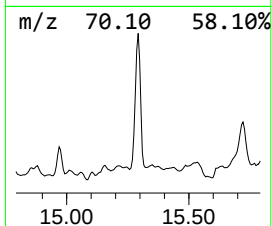
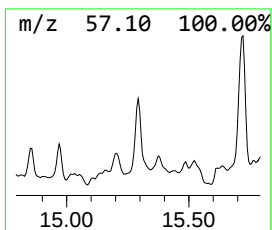
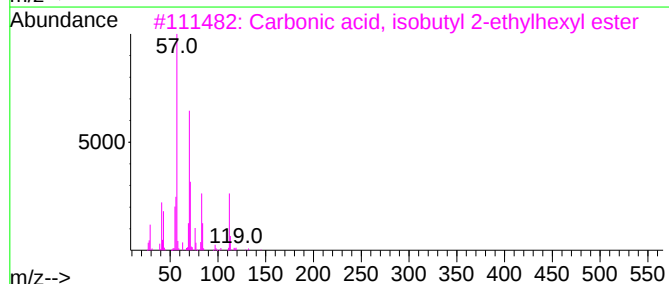
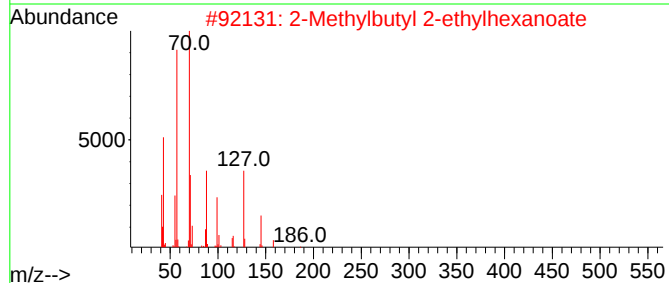
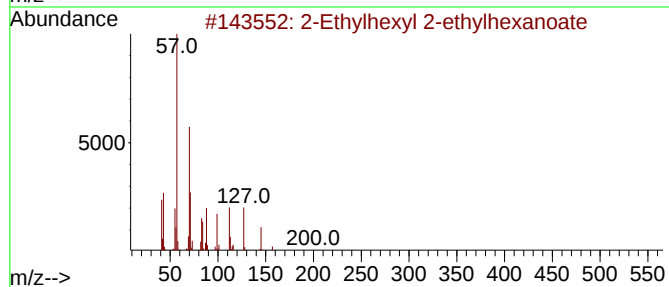
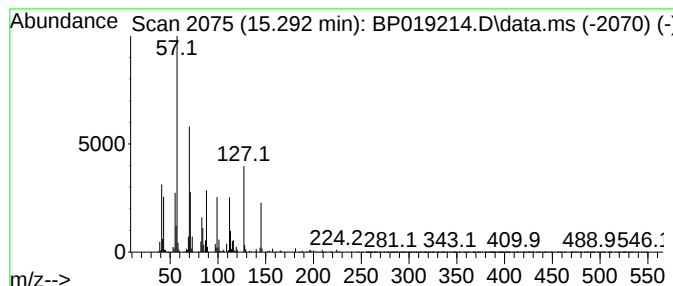
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 2-Ethylhexyl 2-ethylhexanoate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.292	43.40 ng/ul	6979110	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Ethylhexyl 2-ethylhexanoate	256	C16H32O2	007425-14-1	74
2			2-Methylbutyl 2-ethylhexanoate	214	C13H26O2	1010099-98-8	46
3			Carbonic acid, isobutyl 2-ethylh...	230	C13H26O3	1000383-13-3	43
4			1-Ethylpropyl 2-ethylhexanoate	214	C13H26O2	1000099-98-7	43
5			Undecane, 5,5-dimethyl-	184	C13H28	017312-73-1	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
Data File : BP019214.D
Acq On : 02 Jan 2024 18:37
Operator : MA/JU
Sample : 05919-15
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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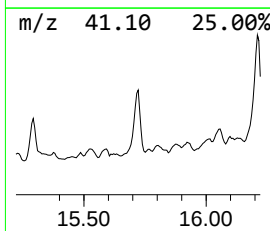
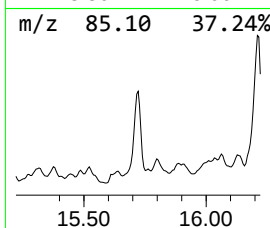
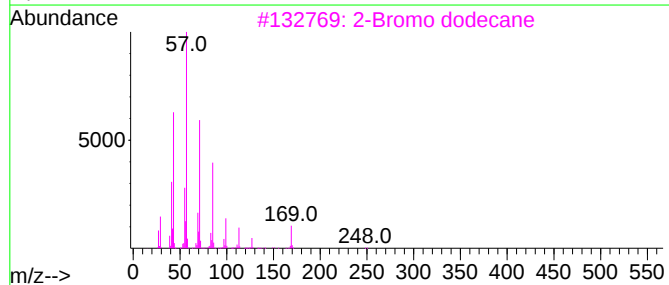
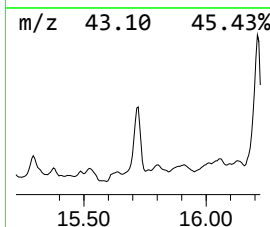
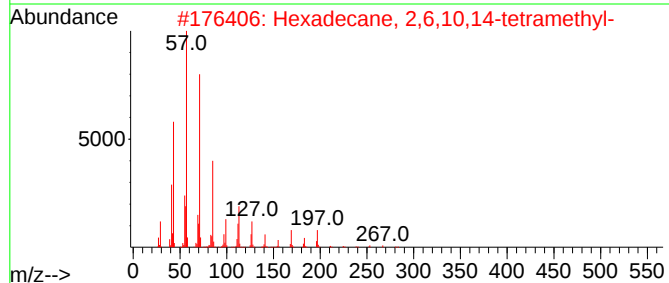
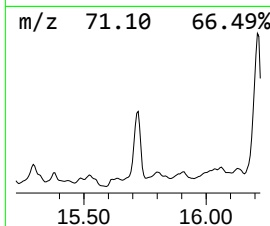
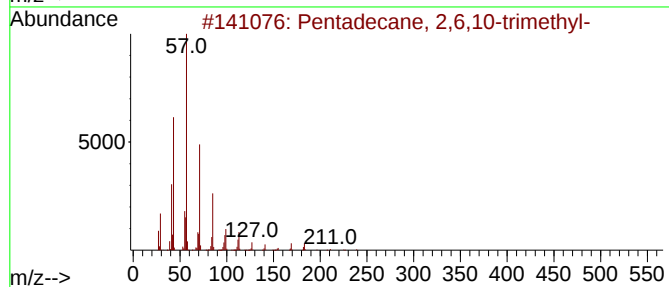
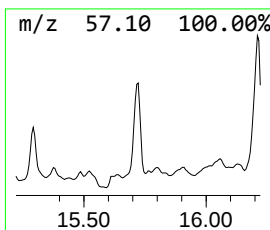
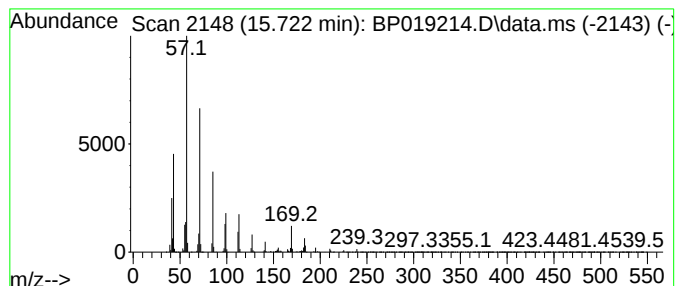
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.722	60.73 ng/ul	9766740	Acenaphthene-d10	14.445

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	94
2		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	80
3		2-Bromo dodecane	248	C12H25Br	013187-99-0	74
4		Hentriacontane	437	C31H64	000630-04-6	72
5		Decane, 3,6-dimethyl-	170	C12H26	017312-53-7	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
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Acq On : 02 Jan 2024 18:37
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Misc :
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Instrument :
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ClientSampleId :
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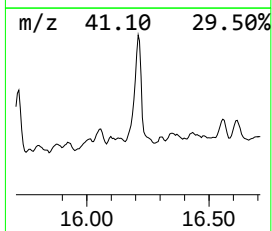
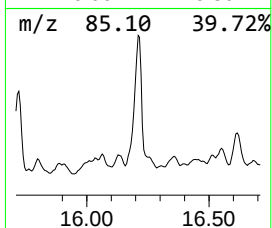
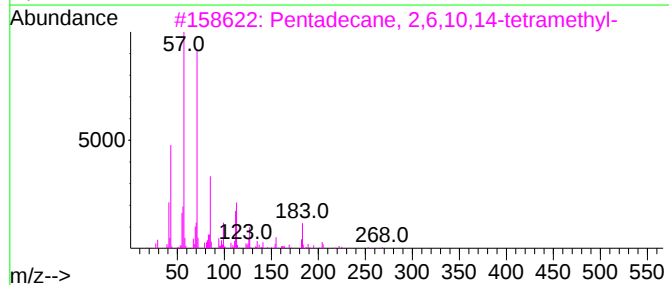
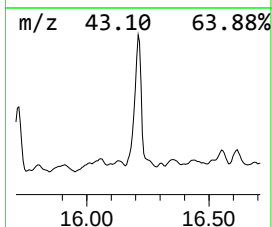
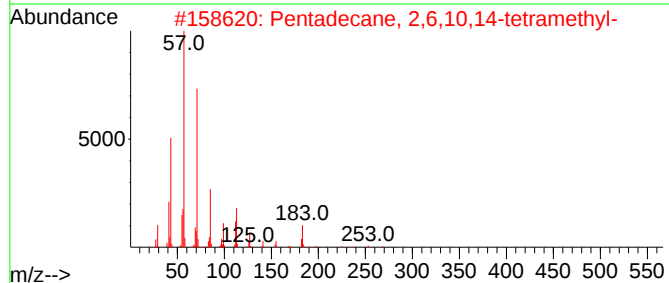
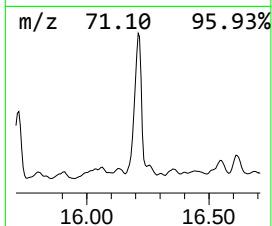
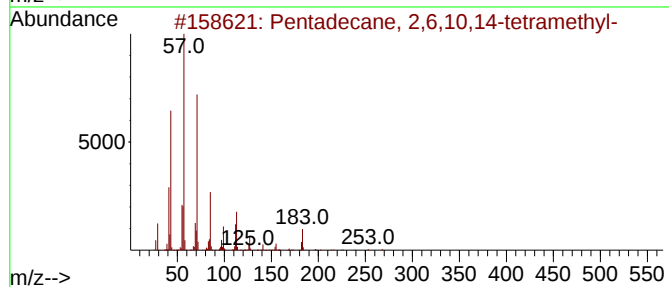
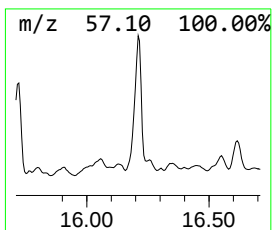
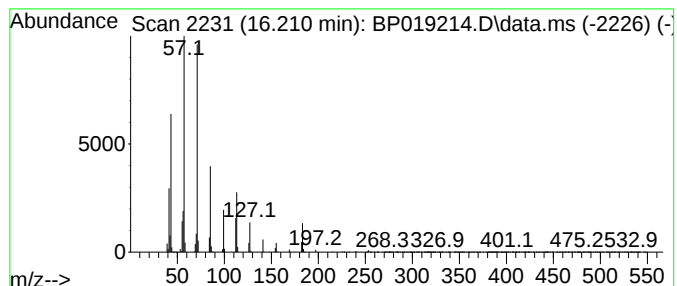
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.210	6.95 ng/ul	16989700	Phenanthrene-d10	17.275

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	90
2			Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	87
3			Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	86
4			Nonane, 2-methyl-5-propyl-	184	C13H28	031081-17-1	53
5			Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
Data File : BP019214.D
Acq On : 02 Jan 2024 18:37
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Sample : 05919-15
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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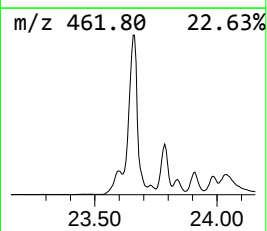
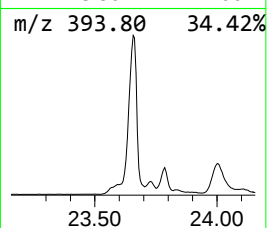
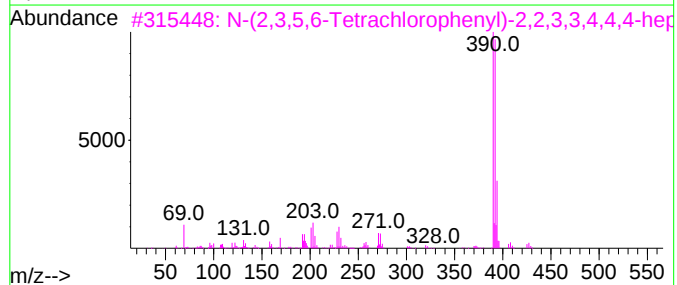
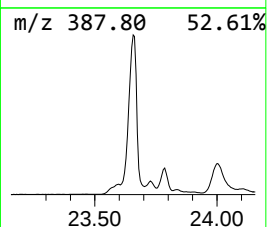
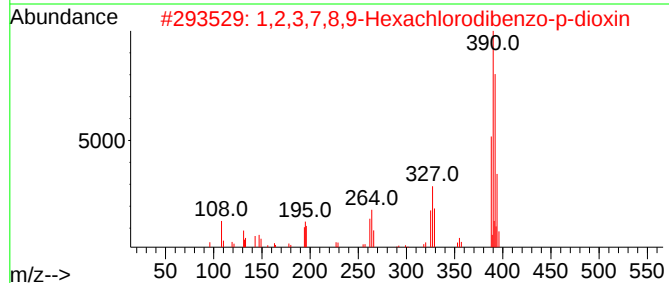
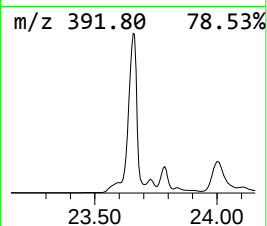
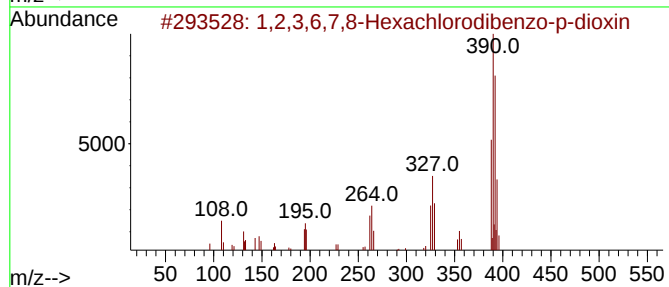
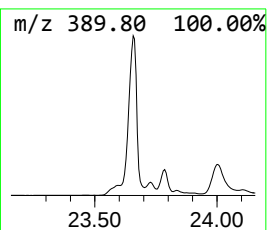
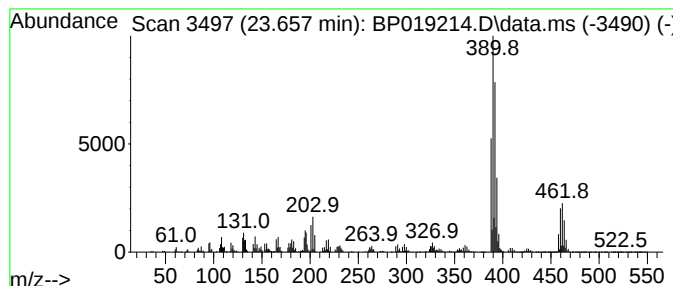
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 1,2,3,6,7,8-Hexachlorodiben... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.657	20.00 ng/ul	30349400	Perylene-d12	23.709

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,3,6,7,8-Hexachlorodibenzo-p-...	388	C12H2Cl6O2	057653-85-7	90		
2	1,2,3,7,8,9-Hexachlorodibenzo-p-...	388	C12H2Cl6O2	019408-74-3	72		
3	N-(2,3,5,6-Tetrachlorophenyl)-2,...	425	C10H2Cl4F7NO	1000373-25-8	40		
4	Thieno[2,3-b]thiophene, 2-bromo-...	390	C11H8Br2N2S2	1000268-03-5	38		
5	4H-1-Benzopyran-2-carboxylic aci...	390	C12H8Br2O5	030113-88-3	30		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
Data File : BP019214.D
Acq On : 02 Jan 2024 18:37
Operator : MA/JU
Sample : 05919-15
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCF47

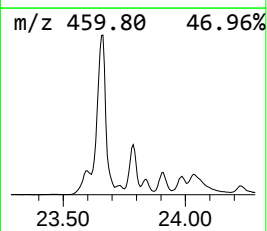
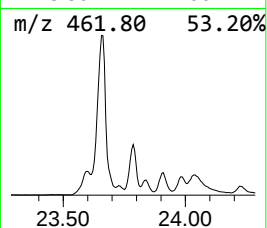
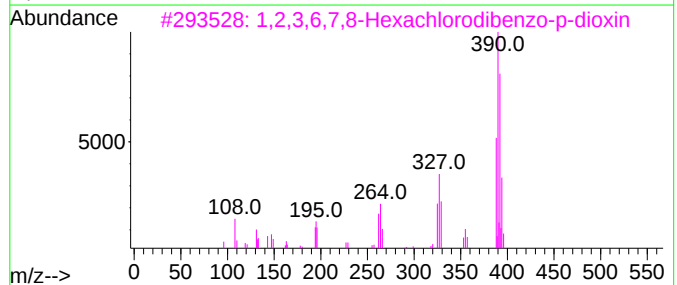
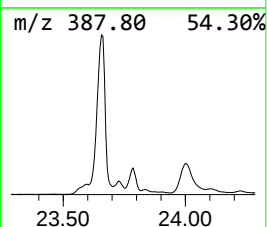
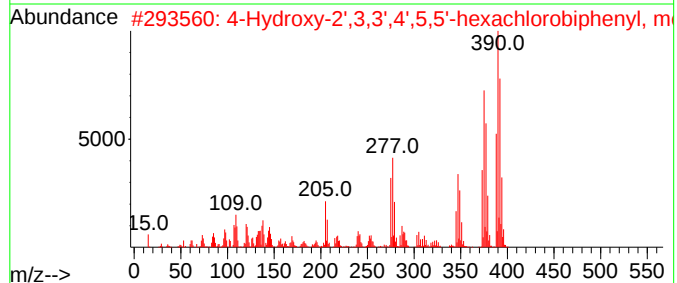
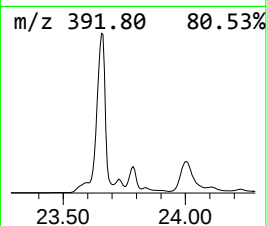
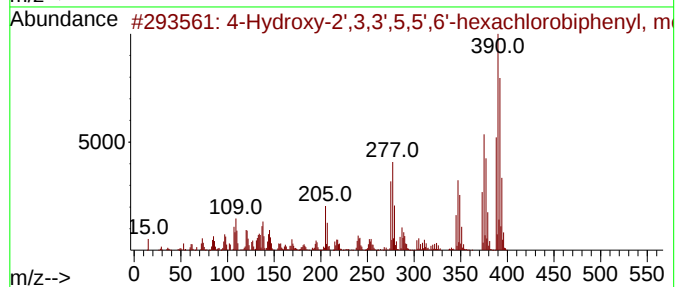
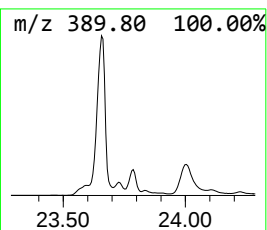
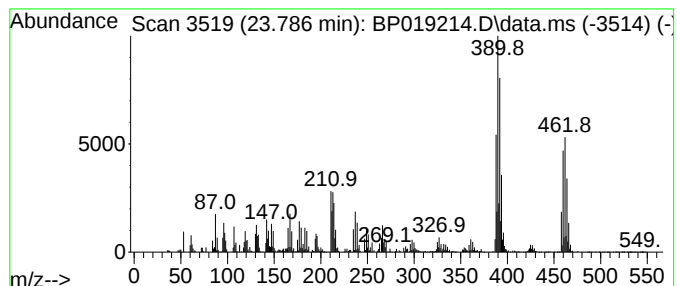
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 4-Hydroxy-2',3,3',5,5',6'-h... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.786	3.50 ng/ul	5313970	Perylene-d12	23.709

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Hydroxy-2',3,3',5,5',6'-hexach...	388	C13H6Cl6O	1000447-88-9	60
2			4-Hydroxy-2',3,3',4',5,5'-hexach...	388	C13H6Cl6O	1000447-88-8	60
3			1,2,3,6,7,8-Hexachlorodibenzo-p-...	388	C12H2Cl6O2	057653-85-7	49
4			N-(2,3,5,6-Tetrachlorophenyl)-2,...	425	C10H2Cl4F7NO	1000373-25-8	45
5			Thieno[2,3-b]thiophene, 2-bromo-...	390	C11H8Br2N2S2	1000268-03-5	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
Data File : BP019214.D
Acq On : 02 Jan 2024 18:37
Operator : MA/JU
Sample : 05919-15
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCF47

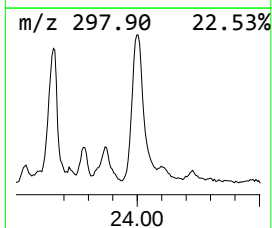
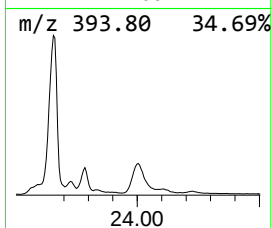
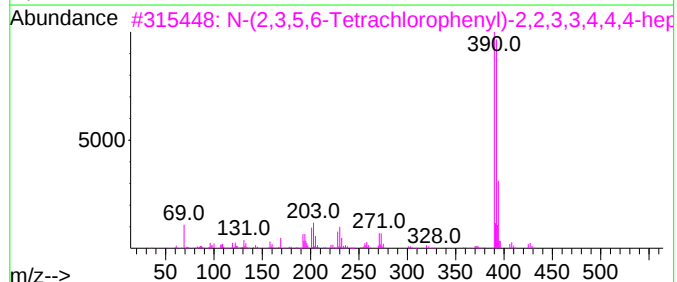
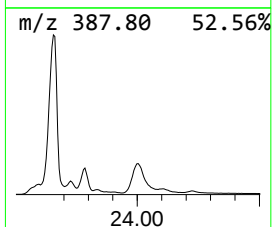
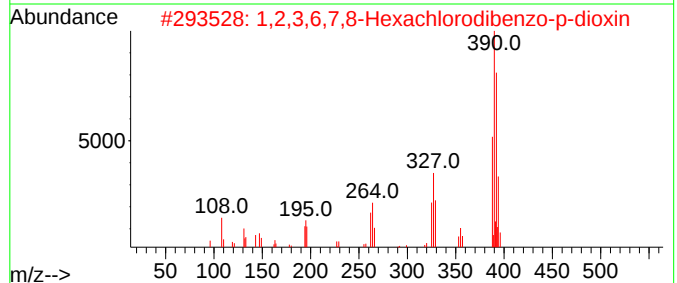
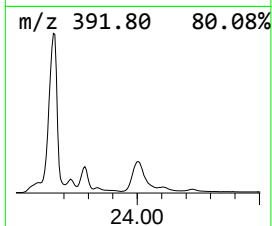
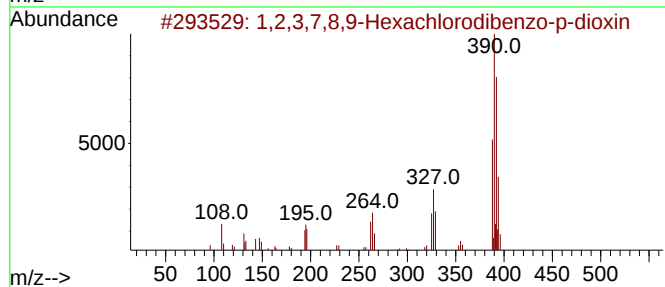
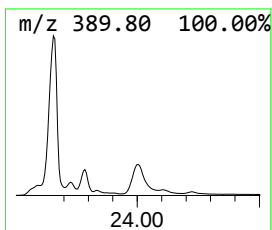
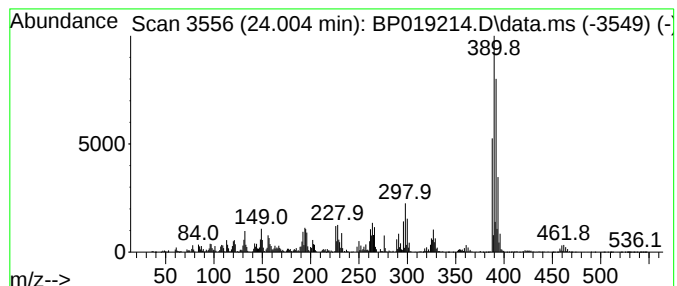
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Quant Title : SVOA CALIBRATION

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

Peak Number 27 1,2,3,7,8,9-Hexachlorodiben... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.004	7.26 ng/ul	11012600	Perylene-d12	23.709

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,3,7,8,9-Hexachlorodibenzo-p-...	388	C12H2Cl6O2	019408-74-3	81		
2	1,2,3,6,7,8-Hexachlorodibenzo-p-...	388	C12H2Cl6O2	057653-85-7	81		
3	N-(2,3,5,6-Tetrachlorophenyl)-2,...	425	C10H2Cl4F7NO	1000373-25-8	50		
4	1-(3-Chlorophenyl)piperazine, N-...	392	C14H12ClF7N2O	1010448-33-8	38		
5	4H-1-Benzopyran-2-carboxylic aci...	390	C12H8Br2O5	030113-88-3	35		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
Data File : BP019214.D
Acq On : 02 Jan 2024 18:37
Operator : MA/JU
Sample : 05919-15
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCF47

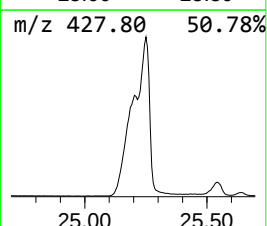
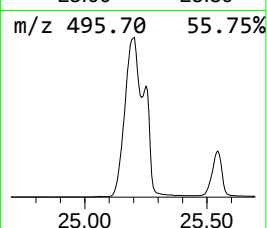
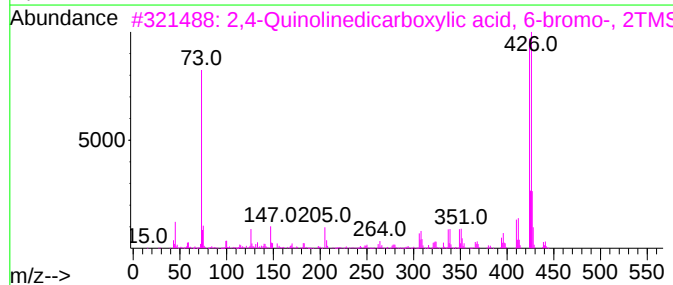
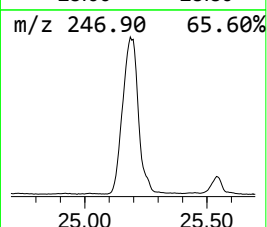
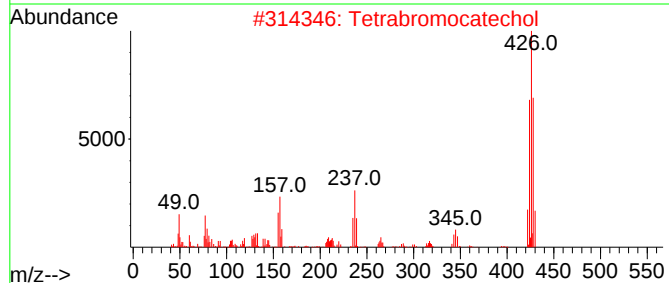
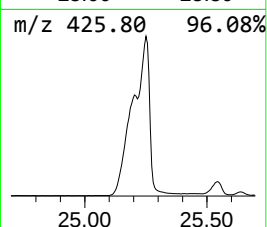
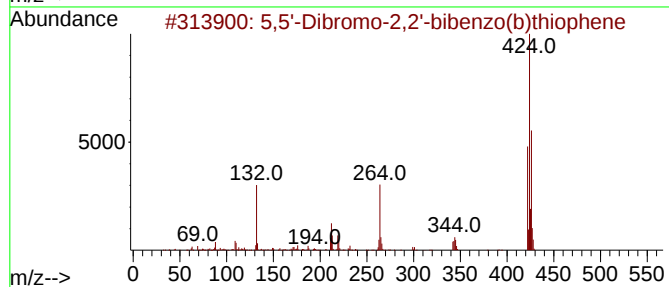
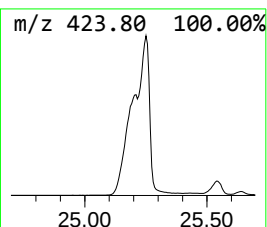
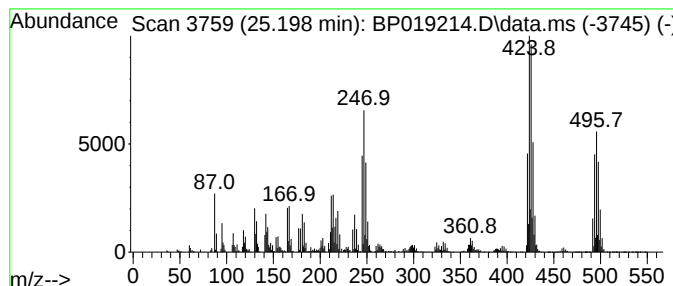
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Quant Title : SVOA CALIBRATION

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

Peak Number 28 unknown-07 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.198	34.21 ng/ul	51906000	Perylene-d12	23.709

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,5'-Dibromo-2,2'-bibenzo(b)thio...	422	C16H8Br2S2	007312-21-2	38		
2	Tetrabromocatechol	422	C6H2Br4O2	000488-47-1	27		
3	2,4-Quinolinedicarboxylic acid, ...	439	C17H22BrN04Si2	1000498-07-4	12		
4	Phenol, 2,3,4,5-tetrabromo-6-met...	420	C7H4Br4O	000576-55-6	12		
5	3,5,3',5'-Tetrakis-trifluorometh...	426	C16H6F12	1000306-81-3	10		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
Data File : BP019214.D
Acq On : 02 Jan 2024 18:37
Operator : MA/JU
Sample : 05919-15
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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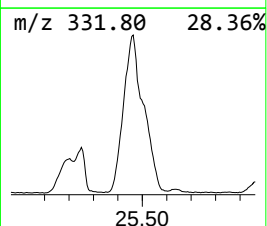
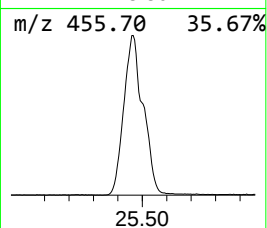
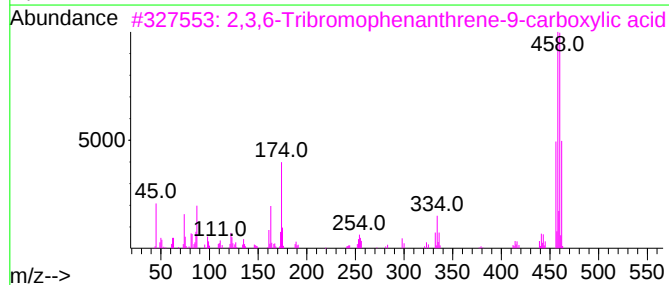
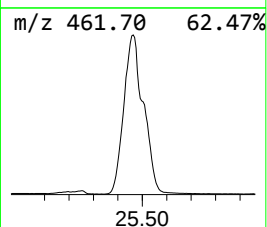
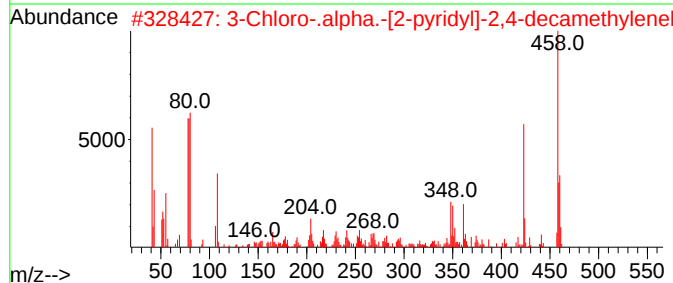
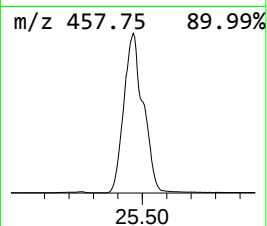
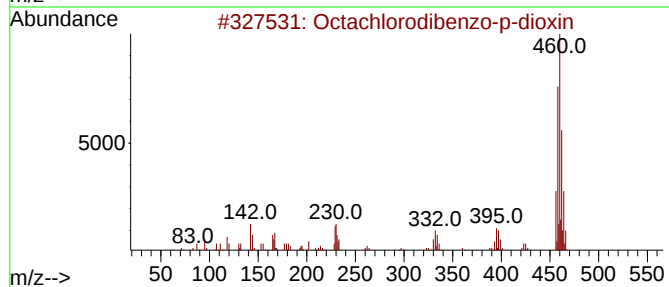
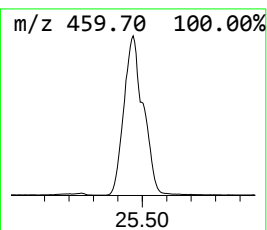
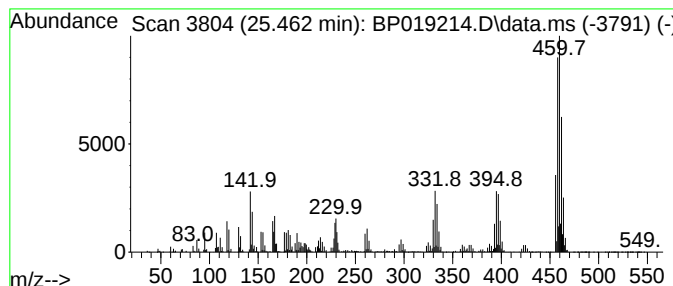
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Quant Title : SVOA CALIBRATION

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

Peak Number 29 Octachlorodibenzo-p-dioxin Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.462	18.23 ng/ul	27657600	Perylene-d12	23.709

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octachlorodibenzo-p-dioxin	456	C12Cl8O2	003268-87-9	91
2			3-Chloro-.alpha.-[2-pyridyl]-2,4...	458	C29H31ClN2O	035158-92-0	60
3			2,3,6-Tribromophenanthrene-9-car...	456	C15H7Br3O2	056988-60-4	37
4			Benzene, 1,4-bis(3-phenylindol-2...	460	C34H24N2	164936-88-3	10
5			Cholest-8(14)-en-3-ol, (3.beta.,...	458	C30H54OSi	055282-50-3	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
Data File : BP019214.D
Acq On : 02 Jan 2024 18:37
Operator : MA/JU
Sample : 05919-15
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCF47

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Hexanol, 2-et...	7.869	3.7	ng/ul	234863	1	7.799	1262610	20.0
Camphor	10.010	4.6	ng/ul	411263	2	10.592	1774170	20.0
(DEL) Alkane: C...	11.269	2.4	ng/ul	212712	2	10.592	1774170	20.0
unknown-01	11.481	2.1	ng/ul	182995	2	10.592	1774170	20.0
(DEL) Alkane: S...	11.622	9.4	ng/ul	834723	2	10.592	1774170	20.0
unknown-02	11.775	2.1	ng/ul	186451	2	10.592	1774170	20.0
(DEL) Alkane: S...	11.869	4.6	ng/ul	409078	2	10.592	1774170	20.0
Hexanoic acid, ...	12.116	2.8	ng/ul	246972	2	10.592	1774170	20.0
unknown-03	12.157	2.2	ng/ul	194419	2	10.592	1774170	20.0
unknown-04	12.328	4.6	ng/ul	407366	2	10.592	1774170	20.0
Phthalic anhydride	12.387	3.4	ng/ul	303239	2	10.592	1774170	20.0
(DEL) Alkane: S...	12.975	15.7	ng/ul	2525890	3	14.445	3216240	20.0
unknown-05	13.022	2.5	ng/ul	395428	3	14.445	3216240	20.0
unknown-06	13.187	3.8	ng/ul	608501	3	14.445	3216240	20.0
(DEL) Alkane: C...	13.245	6.8	ng/ul	1085200	3	14.445	3216240	20.0
(DEL) Alkane: S...	13.357	2.2	ng/ul	347281	3	14.445	3216240	20.0
(DEL) Alkane: S...	13.663	6.0	ng/ul	969938	3	14.445	3216240	20.0
Carbonic acid, ...	13.928	30.8	ng/ul	4949700	3	14.445	3216240	20.0
Carbonic acid, ...	14.357	120.6	ng/ul	19394200	3	14.445	3216240	20.0
2-Bromo dodecane	14.857	7.8	ng/ul	1249080	3	14.445	3216240	20.0
(DEL) Alkane: S...	14.969	24.8	ng/ul	3986350	3	14.445	3216240	20.0
2-Ethylhexyl 2-...	15.292	43.4	ng/ul	6979110	3	14.445	3216240	20.0
(DEL) Alkane: S...	15.722	60.7	ng/ul	9766740	3	14.445	3216240	20.0
(DEL) Alkane: S...	16.210	7.0	ng/ul	16989700	4	17.275	48907000	20.0
1,2,3,6,7,8-Hex...	23.657	20.0	ng/ul	30349400	6	23.709	30349400	20.0
4-Hydroxy-2',3,...	23.786	3.5	ng/ul	5313970	6	23.709	30349400	20.0
1,2,3,7,8,9-Hex...	24.004	7.3	ng/ul	11012600	6	23.709	30349400	20.0
unknown-07	25.198	34.2	ng/ul	51906000	6	23.709	30349400	20.0
Octachlorodiben...	25.462	18.2	ng/ul	27657600	6	23.709	30349400	20.0