

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF053122\
 Data File : BF128602.D
 Acq On : 31 May 2022 18:56
 Operator : CG\JU
 Sample : N3109-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BB-11

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:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052622.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF128602.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.628	55	58	64	rVB2	21183	29584	1.26%	0.113%
2	5.057	468	471	474	rVB	33672	34114	1.46%	0.130%
3	5.498	541	546	555	rBV	2315751	2329843	99.54%	8.904%
4	5.610	560	565	568	rBV	158494	148328	6.34%	0.567%
5	6.522	714	720	729	rBV	2011639	2243633	95.86%	8.575%
6	6.587	729	731	735	rVB	51297	41283	1.76%	0.158%
7	6.840	771	774	778	rBV	717174	641570	27.41%	2.452%
8	7.410	865	871	874	rBV	1770456	1597468	68.25%	6.105%
9	8.128	989	993	996	rBV	956904	782993	33.45%	2.993%
10	9.204	1171	1176	1179	rBV	2906890	2340599	100.00%	8.945%
11	9.434	1206	1215	1223	rVB8	40262	143356	6.12%	0.548%
12	9.881	1288	1291	1295	rVB	938617	783488	33.47%	2.994%
13	10.104	1324	1329	1335	rBV3	35563	69775	2.98%	0.267%
14	10.433	1382	1385	1390	rVB5	30378	37484	1.60%	0.143%
15	10.681	1422	1427	1430	rBV	1534612	1399480	59.79%	5.349%
16	10.780	1439	1444	1454	rVB2	116964	309444	13.22%	1.183%
17	10.992	1477	1480	1481	rBV2	50617	56722	2.42%	0.217%
18	11.010	1481	1483	1492	rVB3	105755	194332	8.30%	0.743%
19	11.369	1541	1544	1547	rBV	774377	718878	30.71%	2.747%
20	11.392	1547	1548	1552	rVB	173603	111388	4.76%	0.426%
21	11.722	1601	1604	1608	rBV4	45347	65656	2.81%	0.251%
22	11.880	1627	1631	1632	rBV3	72141	101202	4.32%	0.387%
23	11.910	1633	1636	1639	rVV4	127382	163789	7.00%	0.626%
24	11.980	1645	1648	1655	rVB3	182221	333951	14.27%	1.276%
25	12.045	1657	1659	1662	rBV3	64089	70991	3.03%	0.271%
26	12.151	1673	1677	1689	rVB4	268924	615819	26.31%	2.354%
27	12.580	1745	1750	1755	rVB2	915797	1315796	56.22%	5.029%
28	12.816	1787	1790	1792	rBV	100539	85084	3.64%	0.325%
29	12.957	1811	1814	1818	rVB	1864026	1764580	75.39%	6.744%
30	13.033	1823	1827	1834	rVB3	203985	442367	18.90%	1.691%
31	13.357	1877	1882	1890	rVB	561291	882031	37.68%	3.371%
32	13.533	1909	1912	1916	rBV2	180214	316241	13.51%	1.209%
33	13.586	1919	1921	1923	rBV	1203250	1395178	59.61%	5.332%
34	14.010	1990	1993	1997	rBV	556392	655714	28.01%	2.506%
35	14.063	1998	2002	2006	rVV2	531080	791446	33.81%	3.025%
36	14.121	2009	2012	2019	rVB2	277113	349323	14.92%	1.335%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052622.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	14.286	2037	2040	2049	rVB	802409	1386732	59.25%	5.300%
38	14.974	2153	2157	2166	rVB	412171	893427	38.17%	3.415%
39	15.486	2241	2244	2253	rVB	388375	522037	22.30%	1.995%

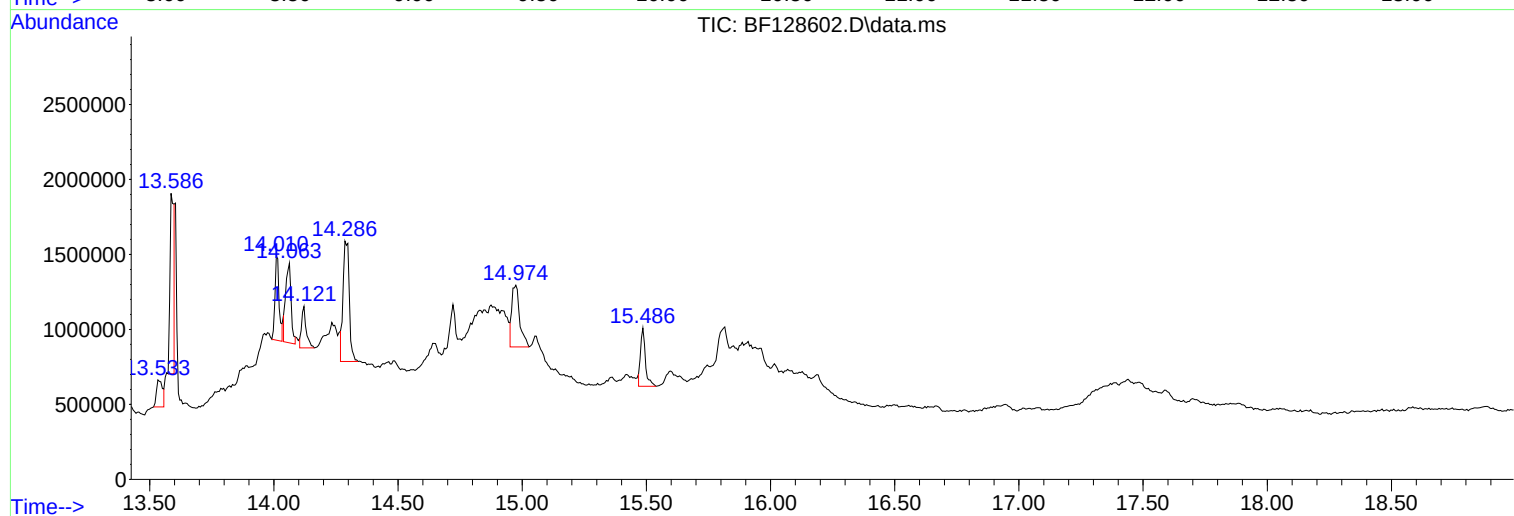
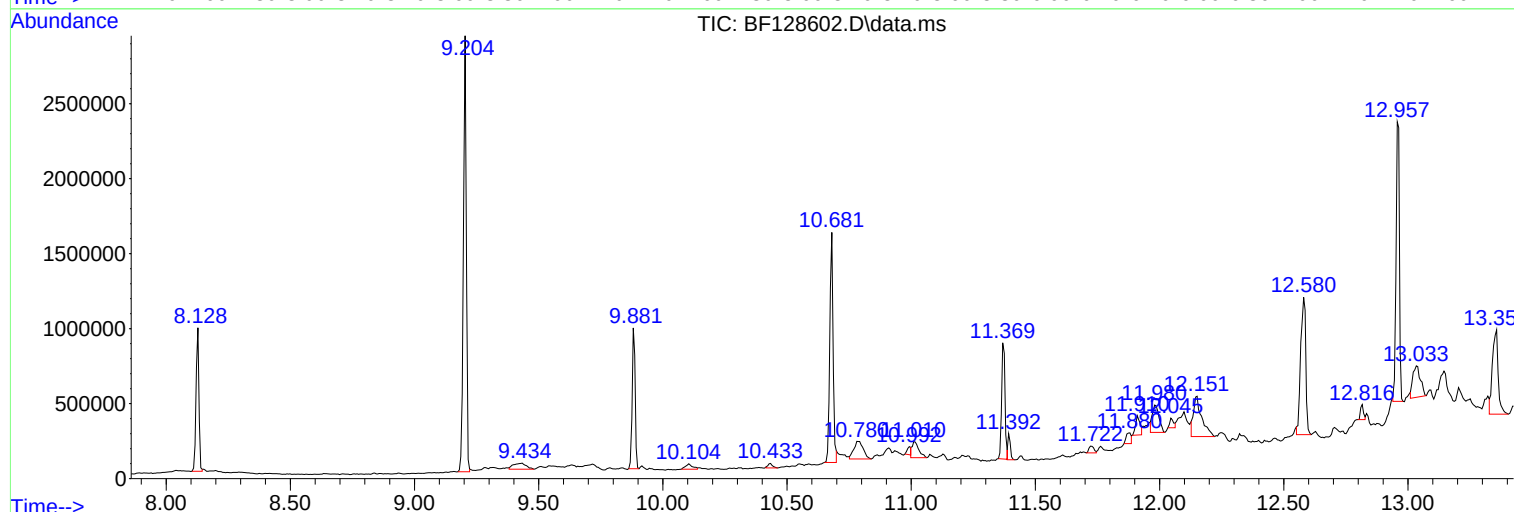
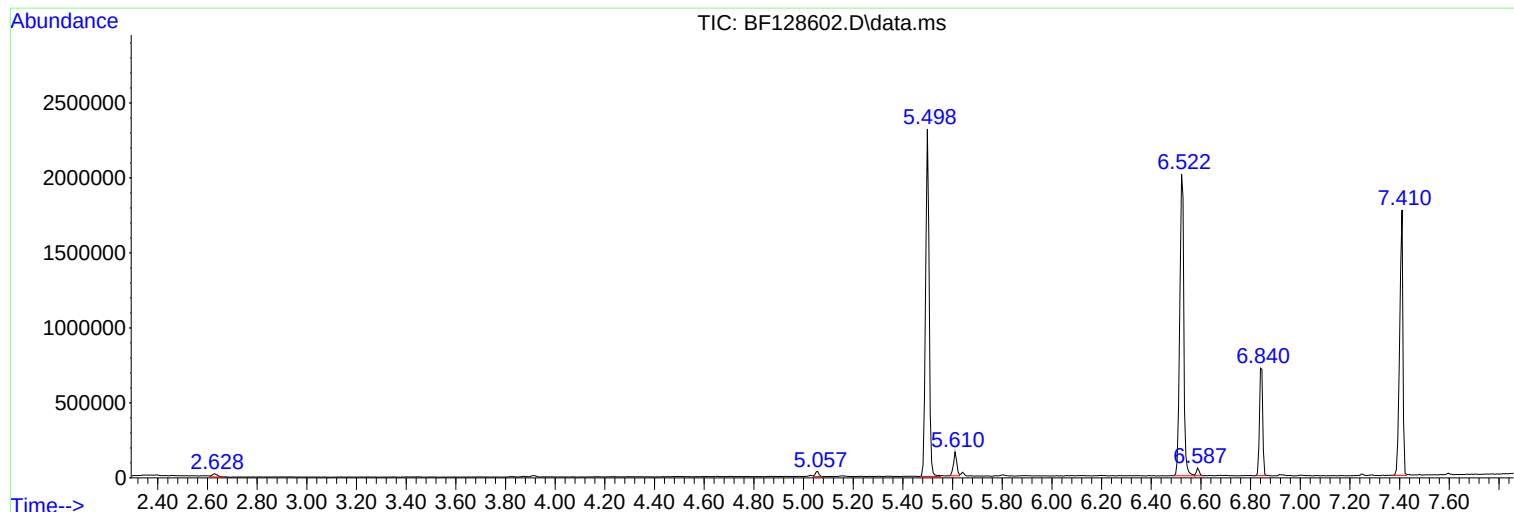
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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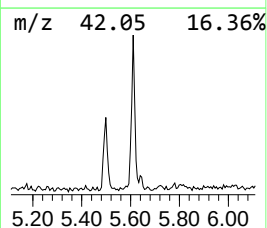
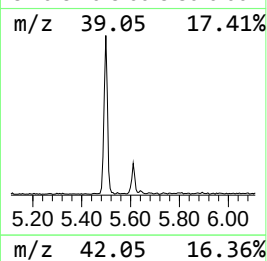
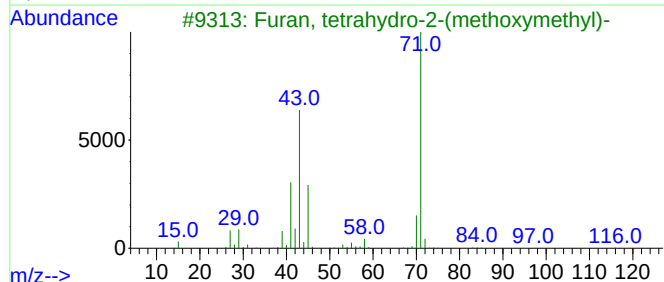
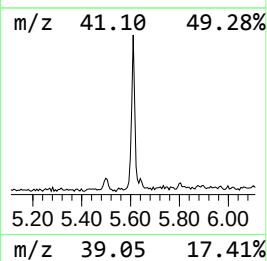
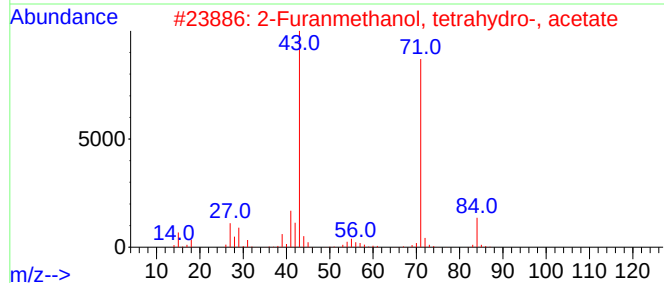
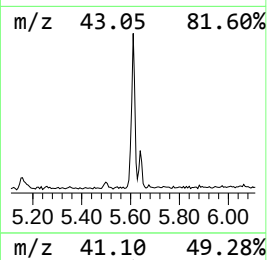
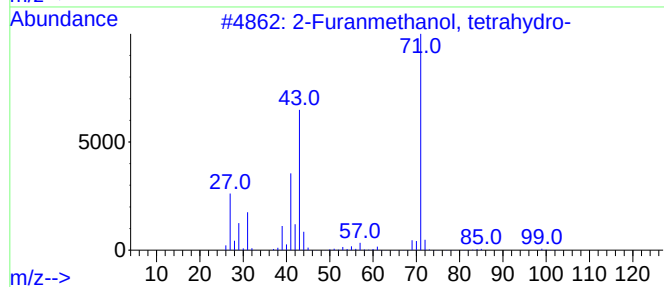
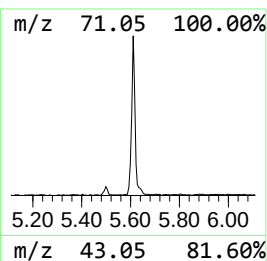
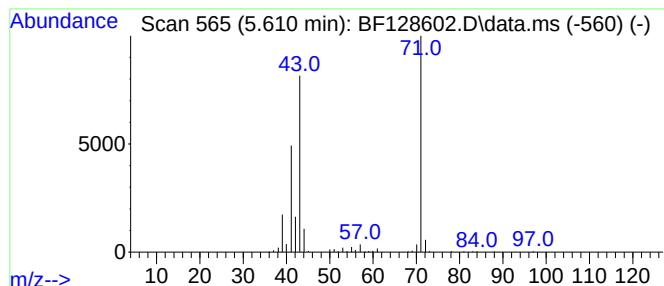
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Furanmethanol, tetrahydro- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.610	4.62 ng	148328	1,4-Dichlorobenzene-d4	6.845

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Furanmethanol, tetrahydro-	102	C5H10O2	000097-99-4	83		
2	2-Furanmethanol, tetrahydro-, ac...	144	C7H12O3	000637-64-9	56		
3	Furan, tetrahydro-2-(methoxymeth...	116	C6H12O2	019354-27-9	56		
4	Butanoyl chloride	106	C4H7ClO	000141-75-3	50		
5	Pentane, 2-bromo-	150	C5H11Br	000107-81-3	40		



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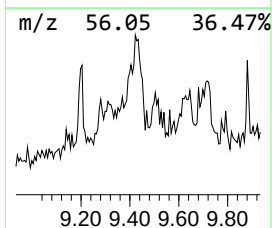
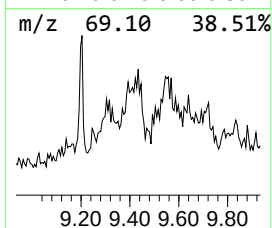
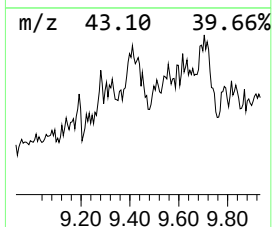
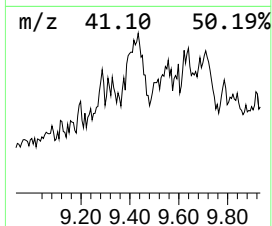
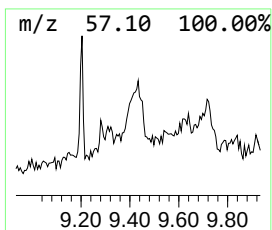
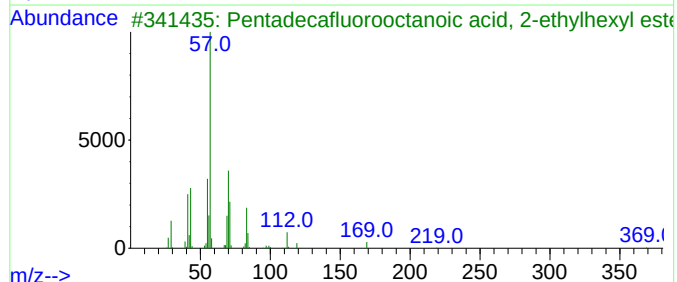
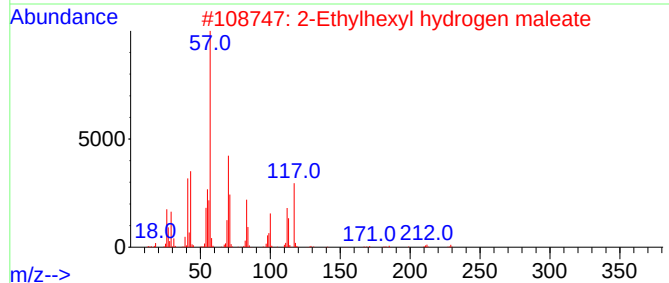
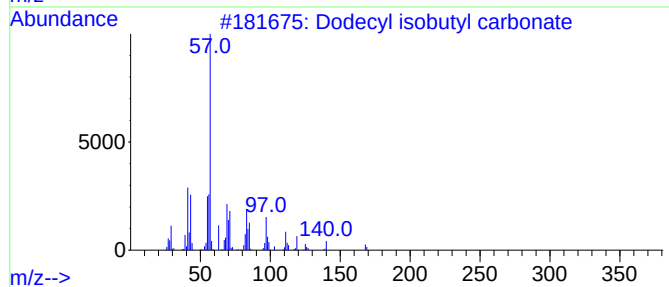
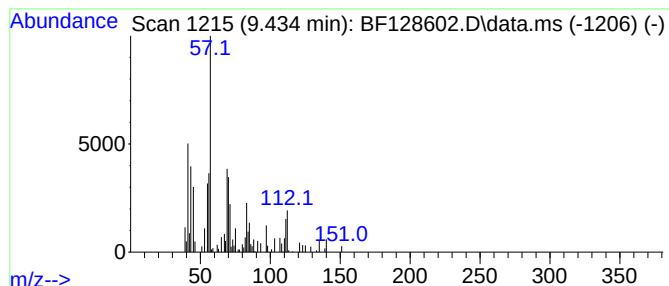
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TIC Library : C:\Database\NIST20.L
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Peak Number 2 Dodecyl isobutyl carbonate Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.434	3.66 ng	143356	Acenaphthene-d10	9.881

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecyl isobutyl carbonate	286	C17H34O3	959067-22-2	59
2			2-Ethylhexyl hydrogen maleate	228	C12H20O4	002370-71-0	50
3			Pentadecafluorooctanoic acid, 2-...	526	C16H17F15O2	1000406-80-9	50
4			3-Octene, (Z)-	112	C8H16	014850-22-7	50
5			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	47



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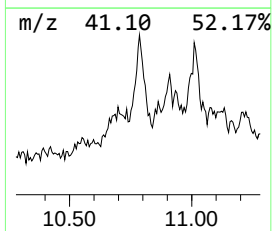
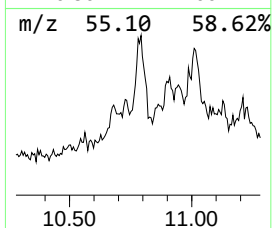
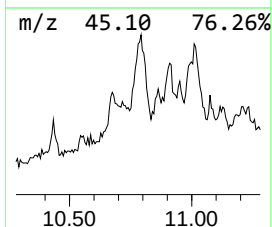
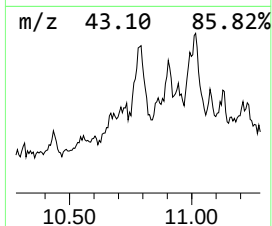
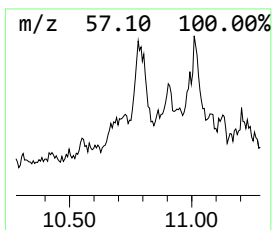
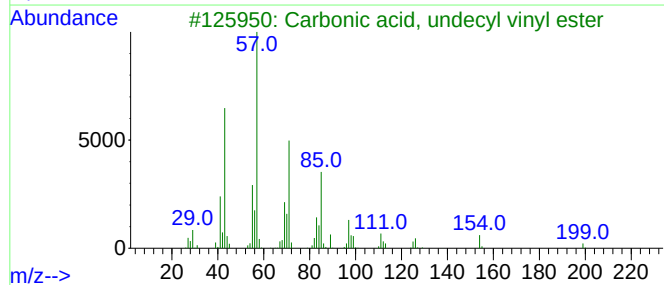
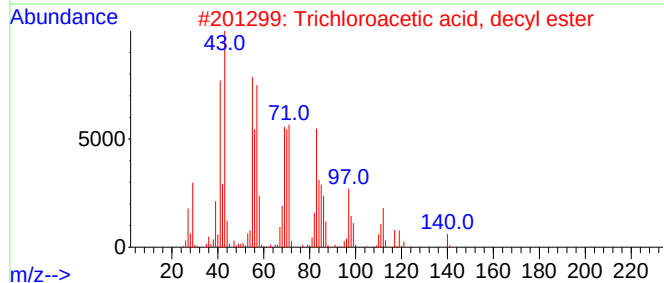
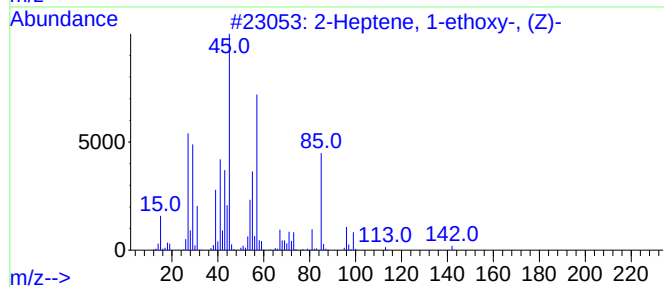
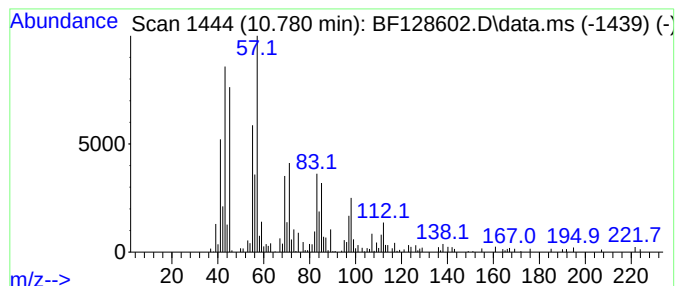
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown10.781 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.781	8.61 ng	309444	Phenanthrene-d10	11.369

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptene, 1-ethoxy-, (Z)-	142	C9H18O	051149-74-7	38		
2	Trichloroacetic acid, decyl ester	302	C12H21Cl3O2	065611-33-8	38		
3	Carbonic acid, undecyl vinyl ester	242	C14H26O3	1000382-54-9	35		
4	Oxalic acid, cyclobutyl tetradec...	340	C20H36O4	1000309-70-9	30		
5	Octatriacontyl pentafluoropropio...	697	C41H77F5O2	1000351-89-1	27		



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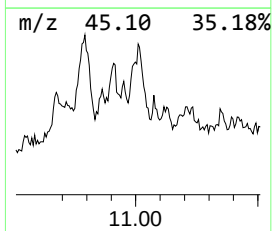
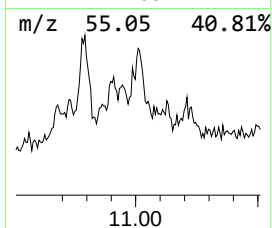
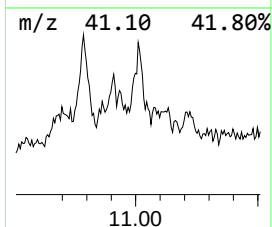
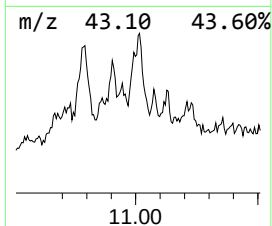
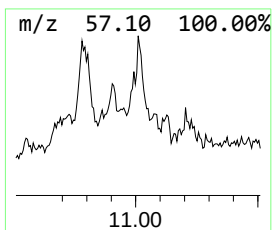
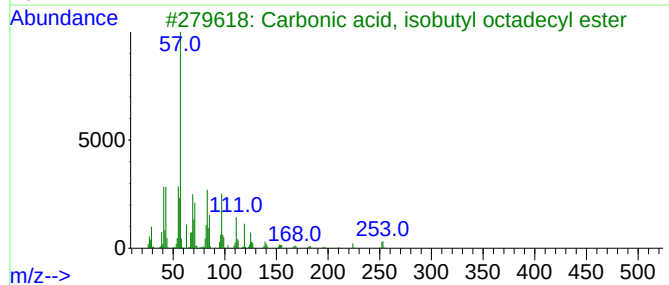
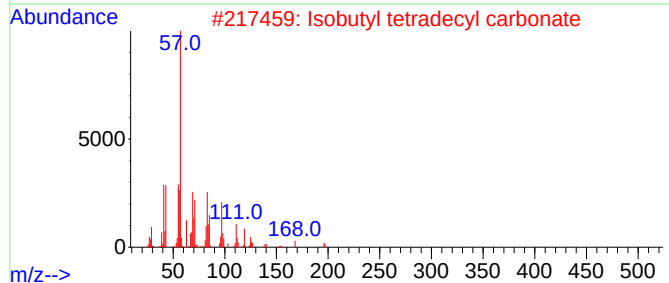
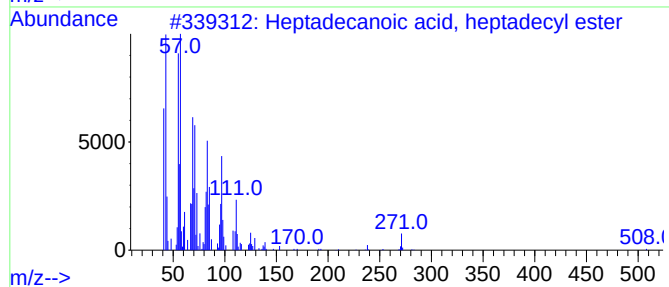
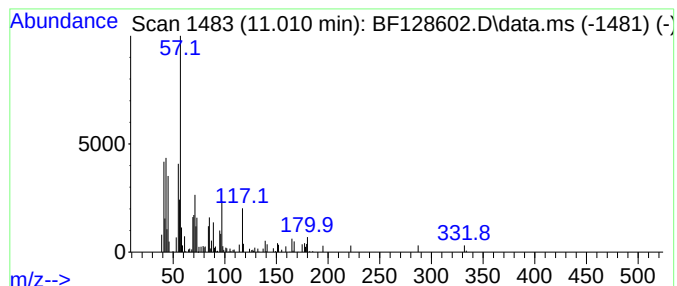
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 unknown11.010 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.010	5.41 ng	194332	Phenanthrene-d10	11.369	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecanoic acid, heptadecyl e...	509	C34H68O2	036617-50-2	47
2	Isobutyl tetradecyl carbonate	314	C19H38O3	959275-58-2	35
3	Carbonic acid, isobutyl octadecy...	370	C23H46O3	1000314-61-5	35
4	Isobutyl tridecyl carbonate	300	C18H36O3	959275-44-6	35
5	Dodecane, 5-cyclohexyl-	252	C18H36	013151-85-4	27



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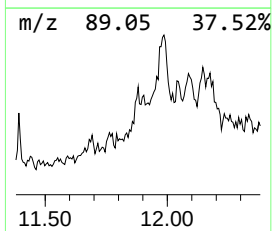
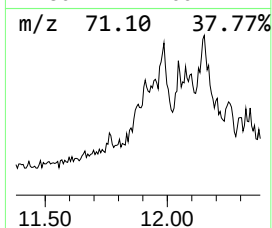
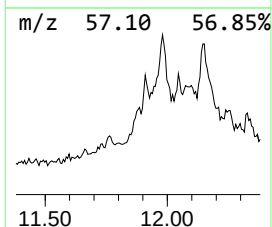
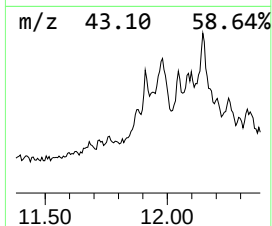
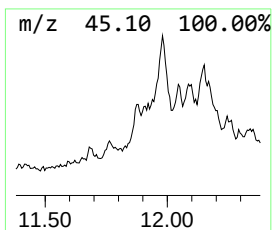
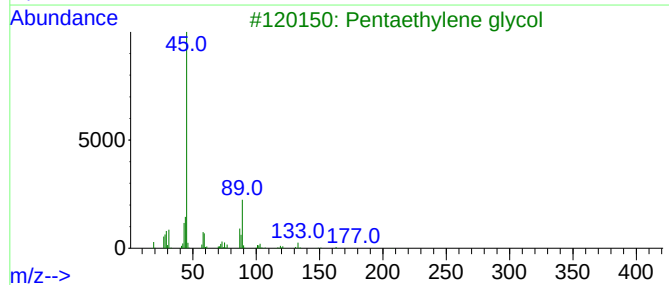
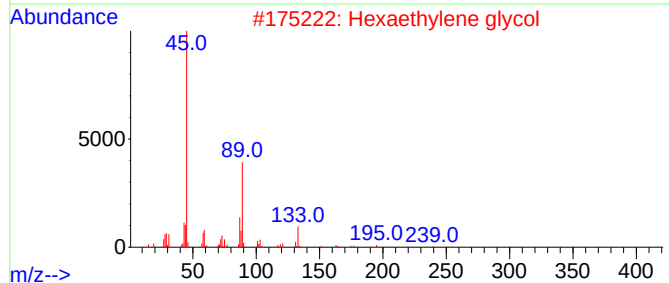
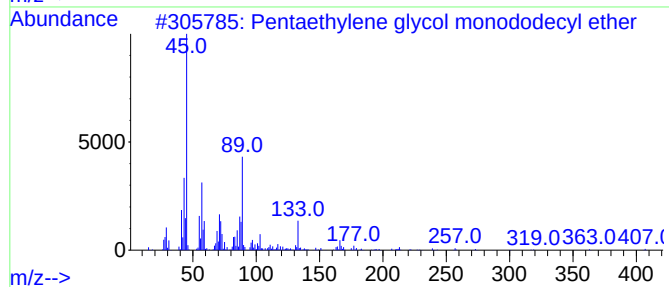
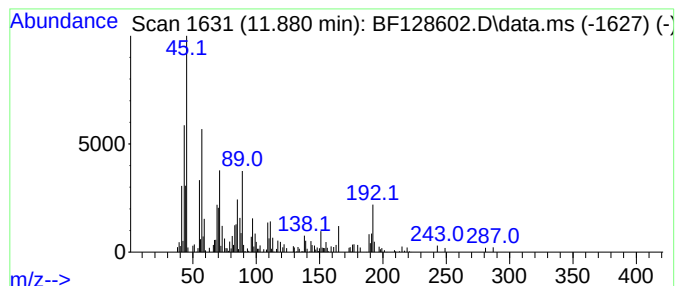
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 unknown11.880 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.880	2.82 ng	101202	Phenanthrene-d10	11.369

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentaethylene glycol monododecyl...	406	C22H46O6	003055-95-6	47
2			Hexaethylene glycol	282	C12H26O7	002615-15-8	46
3			Pentaethylene glycol	238	C10H22O6	004792-15-8	43
4			3,6,9,12-Tetraoxahexadecan-1-ol	250	C12H26O5	001559-34-8	43
5			2-[2-[2-[2-[2-[2-(2-Hydroxyet...	370	C16H34O9	005117-19-1	43



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Acq On : 31 May 2022 18:56
Operator : CG\JU
Sample : N3109-07
Misc :
ALS Vial : 11 Sample Multiplier: 1

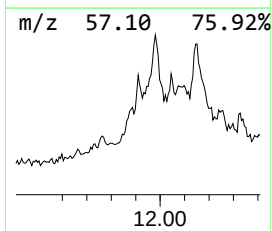
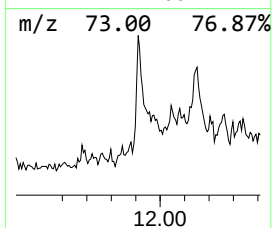
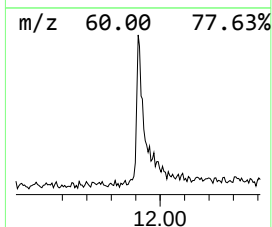
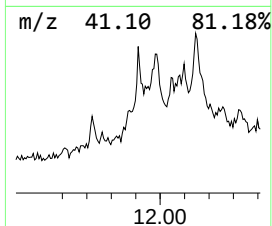
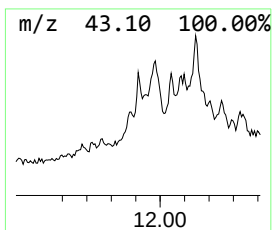
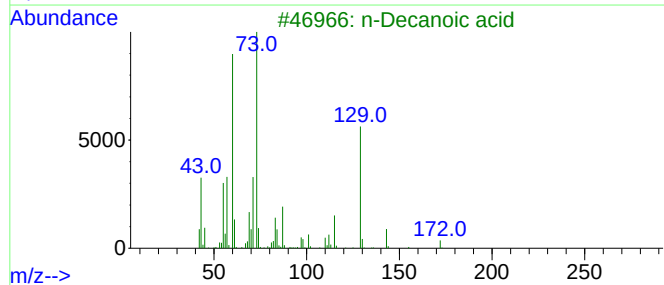
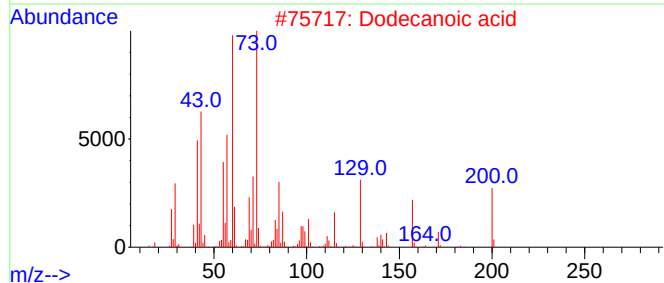
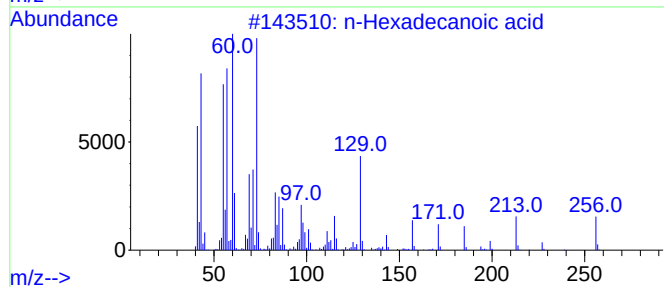
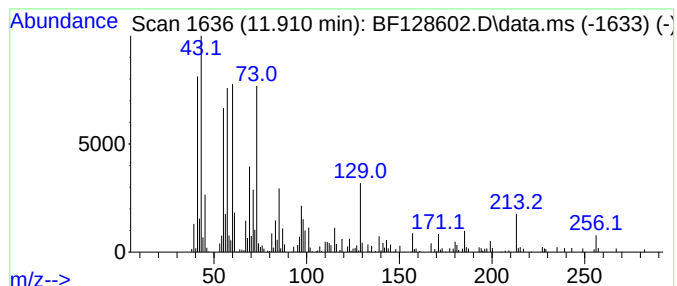
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.910	4.56 ng	163789	Phenanthrene-d10		11.369	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	55
2	Dodecanoic acid		200	C12H24O2	000143-07-7	53
3	n-Decanoic acid		172	C10H20O2	000334-48-5	53
4	8-Methylnonanoic acid		172	C10H20O2	005963-14-4	43
5	Ethanone, 1-(4,5-dihydro-2-thiaz...		129	C5H7NOS	029926-41-8	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF053122\
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Acq On : 31 May 2022 18:56
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ClientSampleId :
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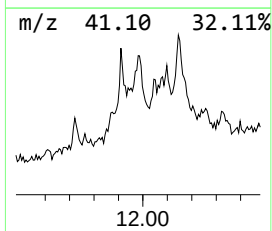
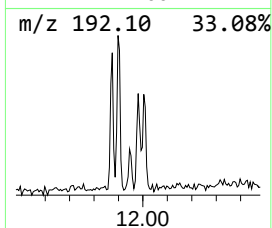
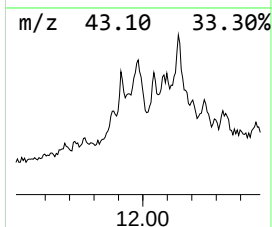
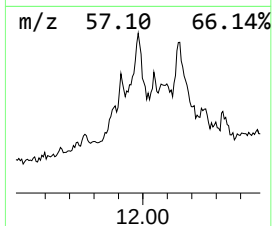
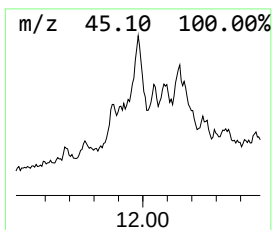
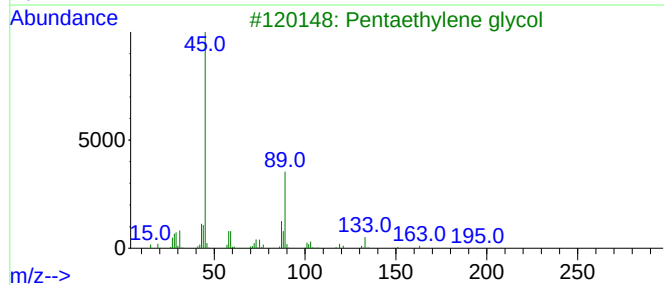
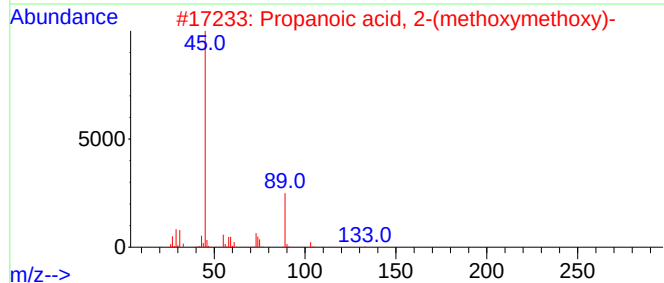
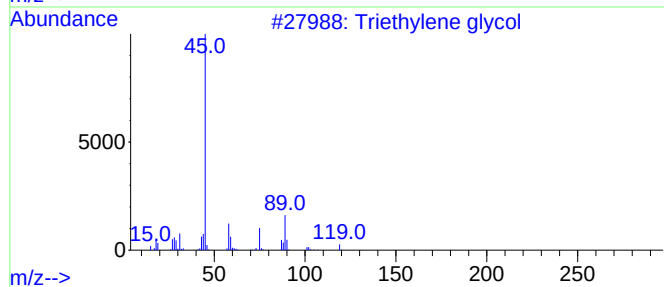
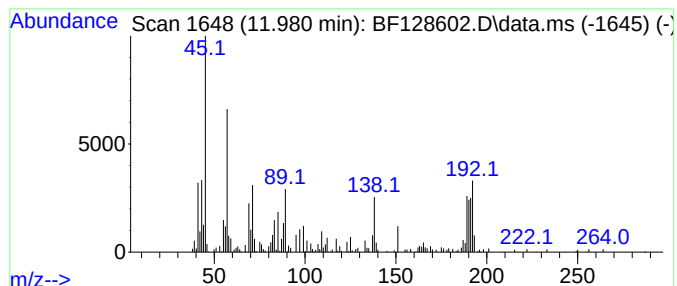
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown11.980 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.980	9.29 ng	333951	Phenanthrene-d10	11.369

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triethylene glycol	150	C6H14O4	000112-27-6	35
2			Propanoic acid, 2-(methoxymethoxy)-	134	C5H10O4	081327-29-9	35
3			Pentaethylene glycol	238	C10H22O6	004792-15-8	35
4			Tetraethylene glycol	194	C8H18O5	000112-60-7	27
5			Diethylene glycol monoethyl ethe...	282	C15H26O3Si	1000376-02-5	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF053122\
Data File : BF128602.D
Acq On : 31 May 2022 18:56
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Sample : N3109-07
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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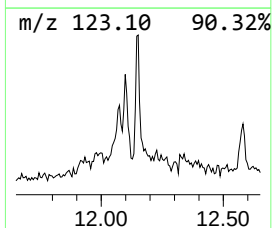
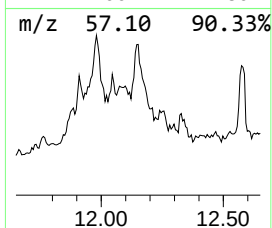
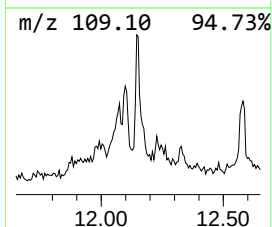
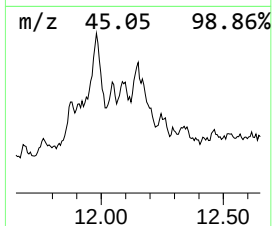
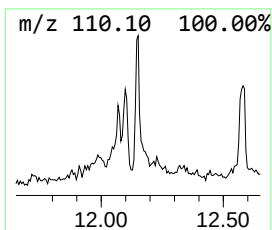
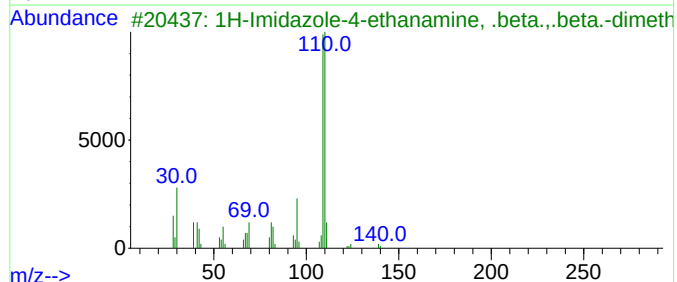
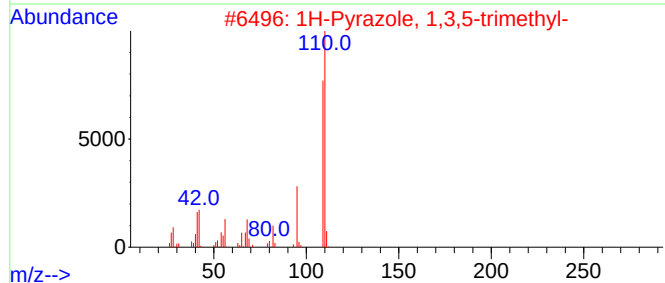
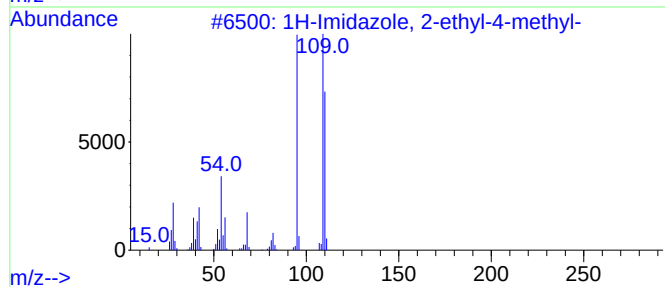
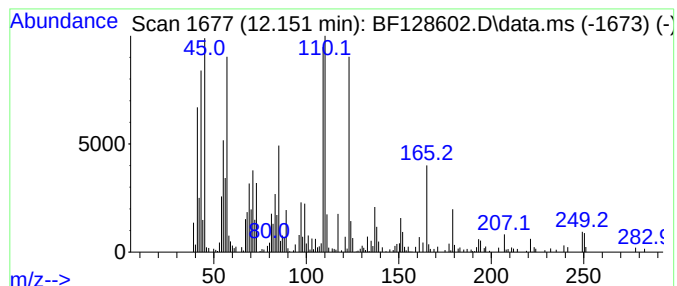
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown12.151 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.151	17.13 ng	615819	Phenanthrene-d10	11.369

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Imidazole, 2-ethyl-4-methyl-	110	C6H10N2	000931-36-2	25
2		1H-Pyrazole, 1,3,5-trimethyl-	110	C6H10N2	001072-91-9	25
3		1H-Imidazole-4-ethanamine, .beta...	139	C7H13N3	021150-01-6	22
4		Naphthalene, decahydro-2,6-dimet...	278	C20H38	054964-85-1	11
5		2,4-Heptadienal, 2,4-dimethyl-	138	C9H14O	042452-48-2	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF053122\
Data File : BF128602.D
Acq On : 31 May 2022 18:56
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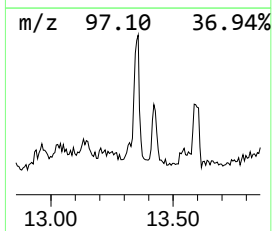
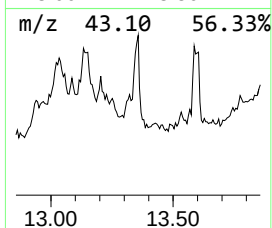
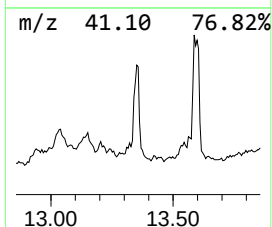
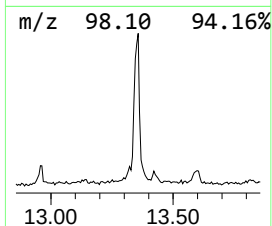
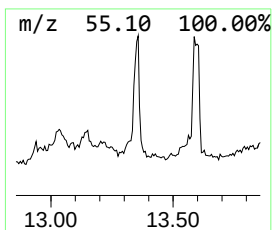
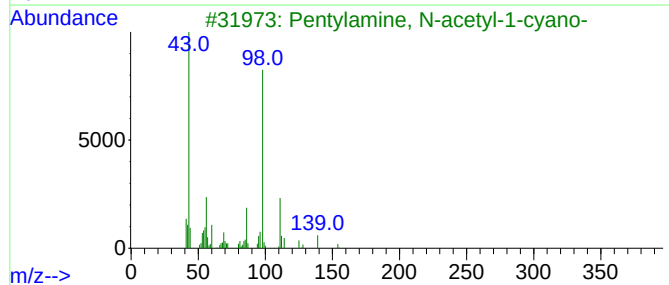
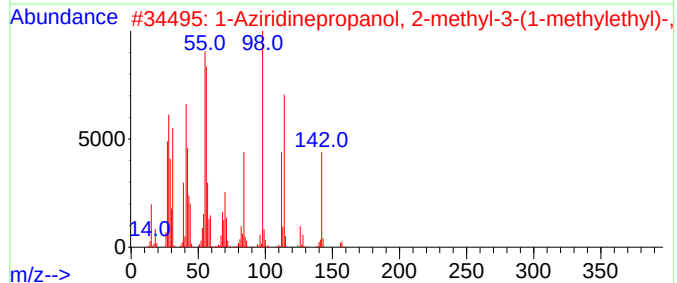
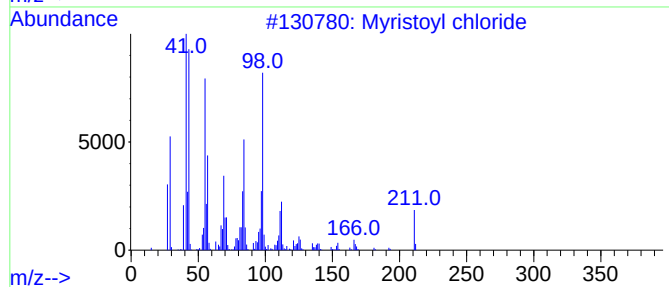
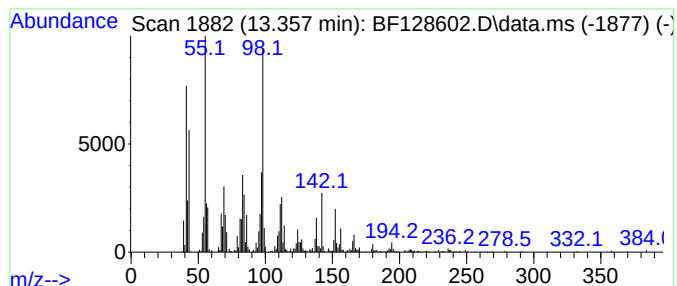
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown13.357 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.357	26.90 ng	882031	Chrysene-d12	14.016

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Myristoyl chloride	246	C14H27ClO	000112-64-1	49
2			1-Aziridinepropanol, 2-methyl-3-...	157	C9H19NO	055669-83-5	38
3			Pentylamine, N-acetyl-1-cyano-	154	C8H14N2O	027395-08-0	38
4			Cyclononanone	140	C9H16O	003350-30-9	35
5			Cyclohexanone, 2-propyl-	140	C9H16O	000094-65-5	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF053122\
Data File : BF128602.D
Acq On : 31 May 2022 18:56
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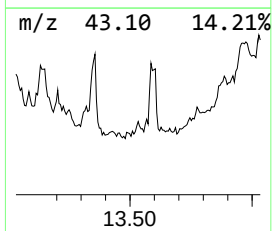
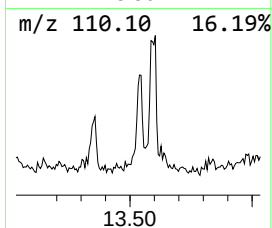
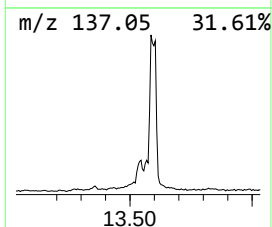
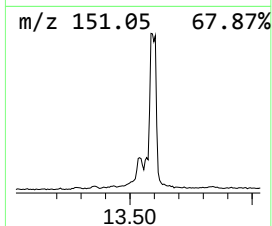
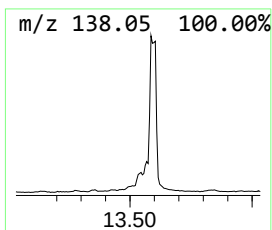
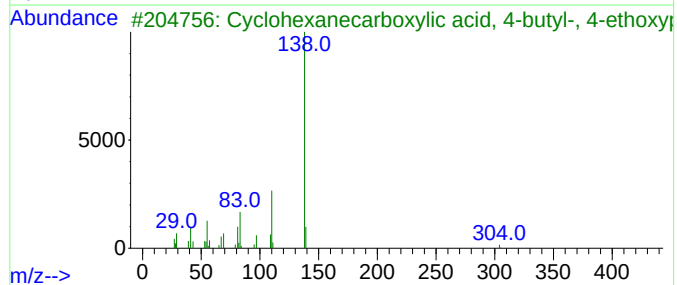
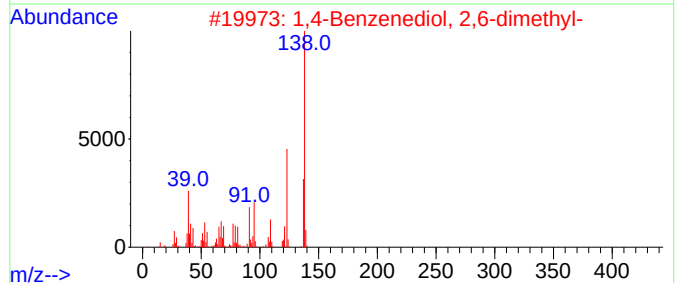
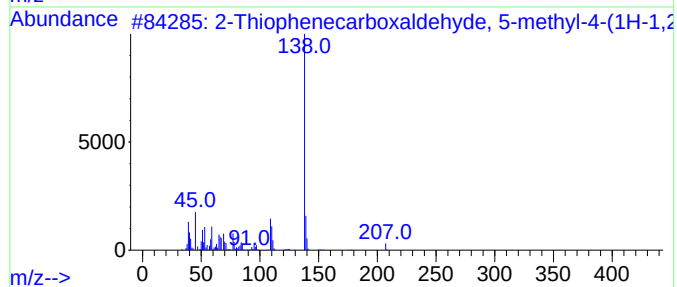
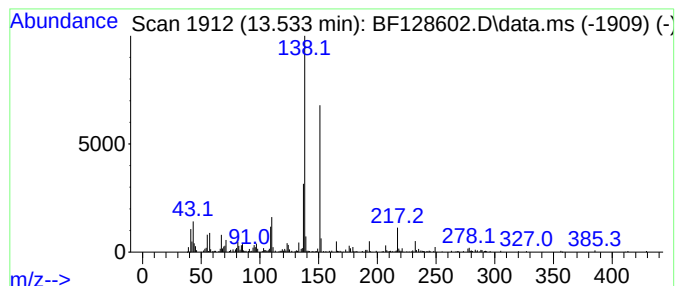
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown13.533 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.533	9.65 ng	316241	Chrysene-d12	14.016

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Thiophenecarboxaldehyde, 5-met...	207	C9H9N3O5	1000337-27-9	46		
2	1,4-Benzenediol, 2,6-dimethyl-	138	C8H10O2	000654-42-2	43		
3	Cyclohexanecarboxylic acid, 4-bu...	304	C19H28O3	067589-47-3	38		
4	3,5-Diaminobenzyl alcohol	138	C7H10N2O	028150-13-2	38		
5	Cyclohexanecarboxylic acid, 4-pr...	290	C18H26O3	067589-39-3	38		



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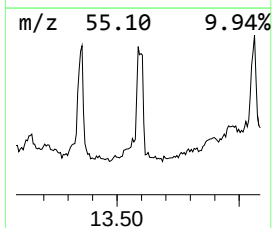
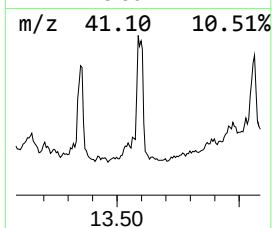
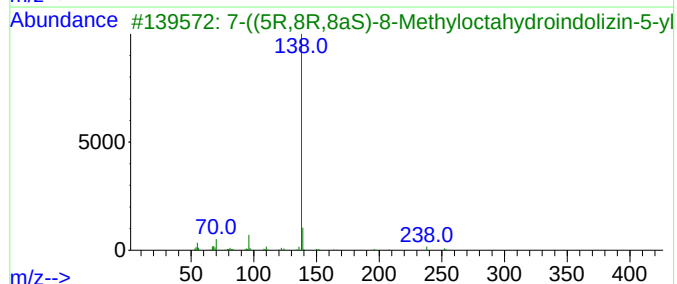
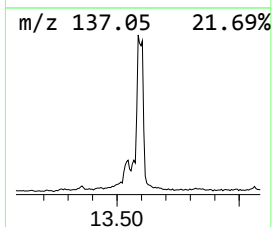
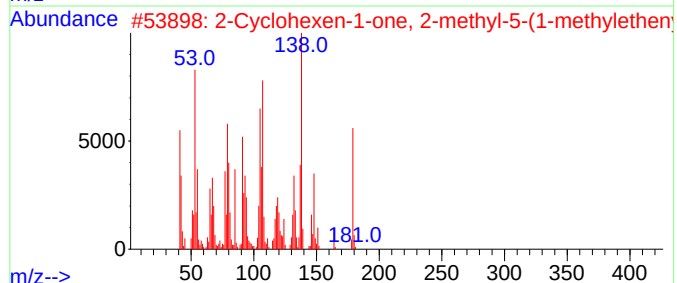
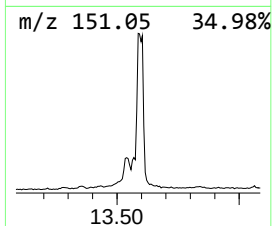
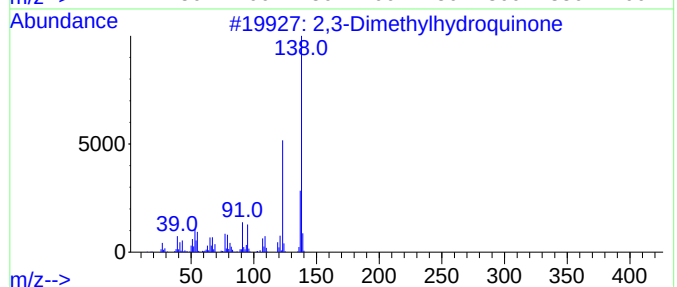
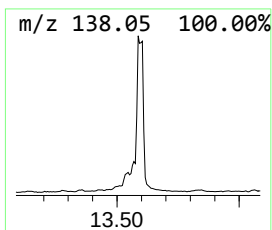
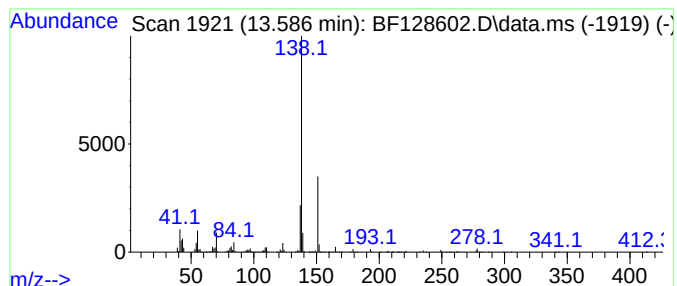
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 2,3-Dimethylhydroquinone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.586	42.55 ng	1395180	Chrysene-d12	14.016

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2,3-Dimethylhydroquinone	138	C8H10O2	000608-43-5	52
2		2-Cyclohexen-1-one, 2-methyl-5-(...	179	C11H17NO	057397-12-3	52
3		7-((5R,8R,8aS)-8-Methyloctahydro...	253	C16H31NO	1000367-28-5	47
4		Gephyrotoxin 235b'	235	C16H29N	120030-10-6	47
5		Gephyrotoxin 207a	207	C14H25N	141724-48-3	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF053122\
Data File : BF128602.D
Acq On : 31 May 2022 18:56
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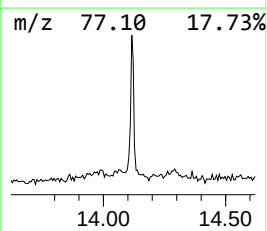
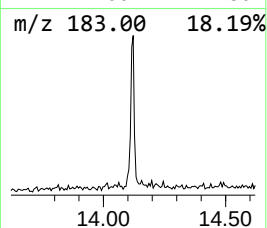
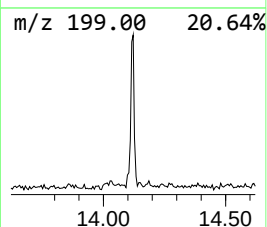
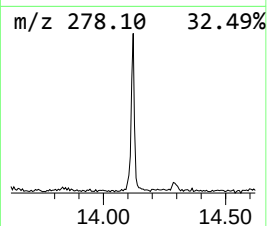
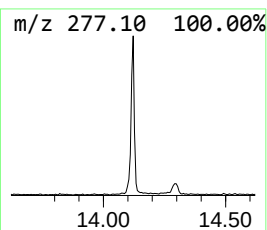
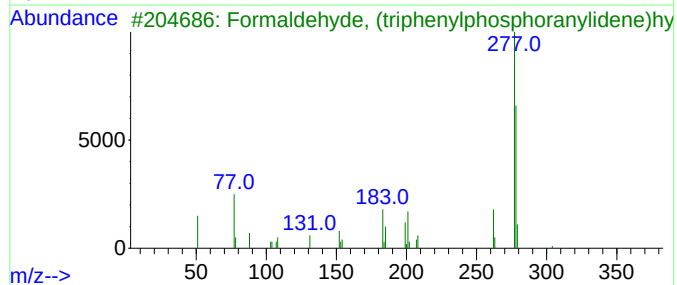
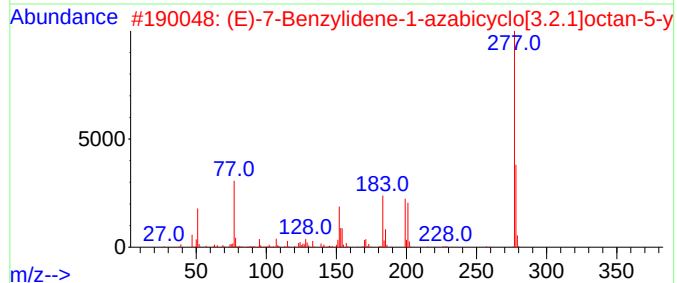
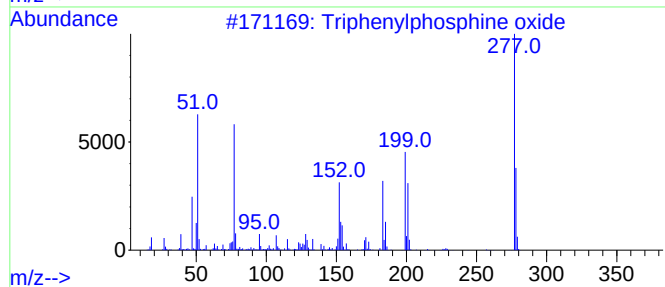
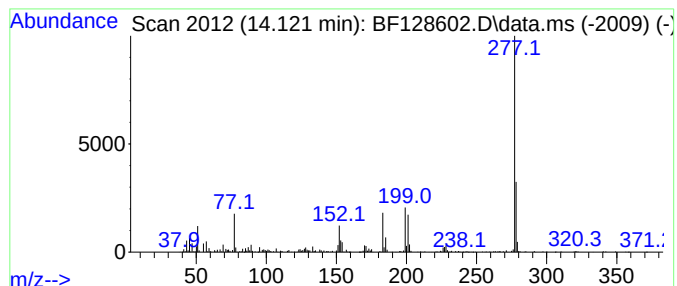
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Triphenylphosphine oxide Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.121	10.65 ng	349323	Chrysene-d12	14.016

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triphenylphosphine oxide	278	C18H15OP	000791-28-6	97
2			(E)-7-Benzylidene-1-azabicyclo[3...]	293	C15H19N03S	1000483-83-4	90
3			Formaldehyde, (triphenylphosphor...	304	C19H17N2P	015990-54-2	45
4			(Carbethoxymethyl)-triphenylphos...	428	C22H22BrO2P	001530-45-6	40
5			3-(3,4-Dichlorophenyl)-2-thioxo-...	277	C9H5Cl2NOS2	017046-34-3	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF053122\
Data File : BF128602.D
Acq On : 31 May 2022 18:56
Operator : CG\JU
Sample : N3109-07
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BB-11

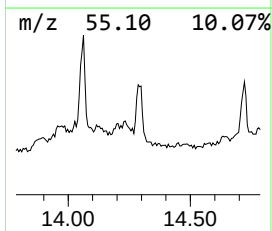
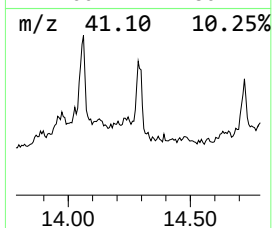
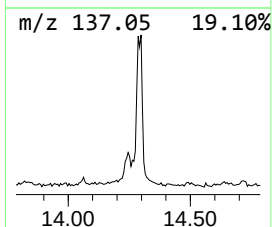
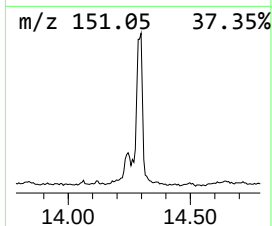
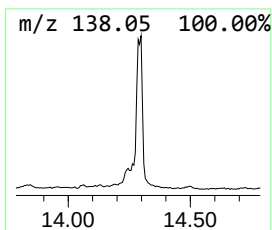
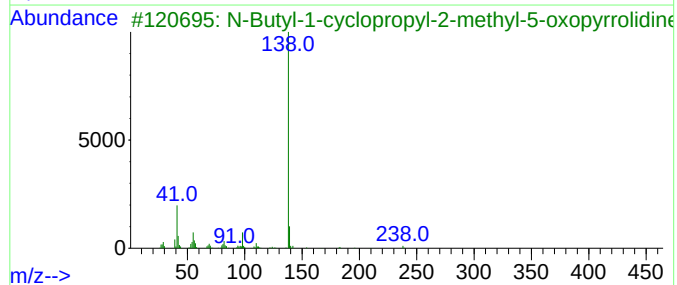
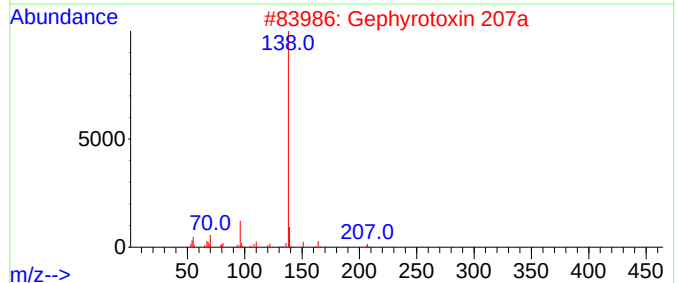
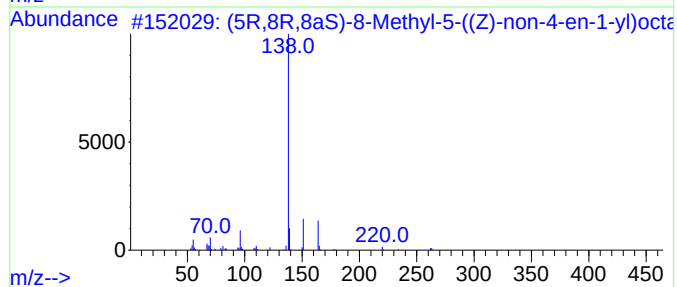
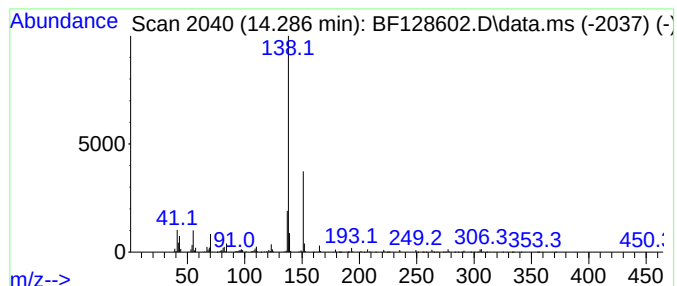
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 (5R,8R,8aS)-8-Methyl-5-((Z)... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.286	42.30 ng	1386730	Chrysene-d12	14.016

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(5R,8R,8aS)-8-Methyl-5-((Z)-non-...	263	C18H33N	1000367-14-0	50
2			Gephyrotoxin 207a	207	C14H25N	141724-48-3	47
3			N-Butyl-1-cyclopropyl-2-methyl-5...	238	C13H22N2O2	1010458-31-0	43
4			2,5-Cyclohexadiene-1,4-dione, 2-...	138	C7H6O3	000615-91-8	43
5			Gephyrotoxin 235b'	235	C16H29N	120030-10-6	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF053122\
Data File : BF128602.D
Acq On : 31 May 2022 18:56
Operator : CG\JU
Sample : N3109-07
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BB-11

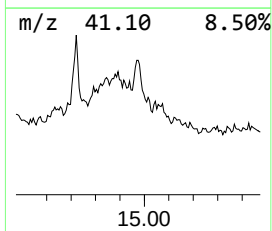
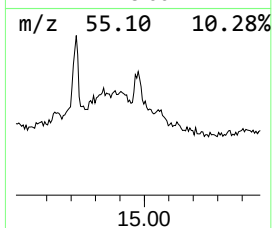
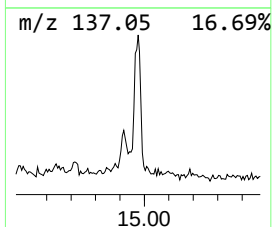
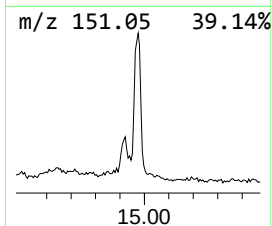
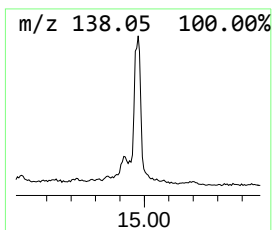
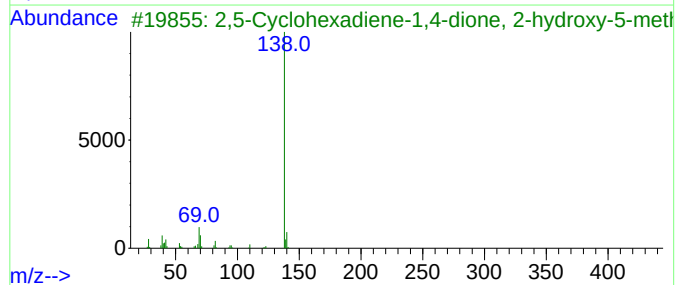
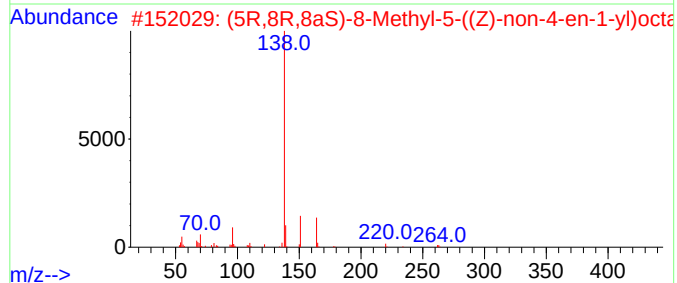
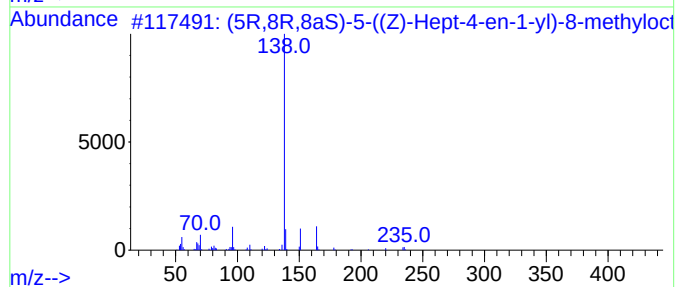
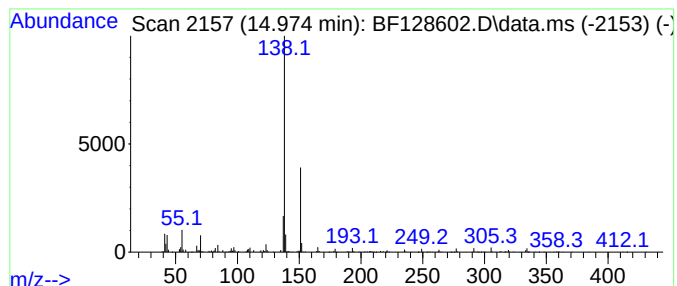
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown14.974 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.974	34.23 ng	893427	Perylene-d12	15.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(5R,8R,8aS)-5-((Z)-Hept-4-en-1-y...	235	C16H29N	109175-47-5	49		
2	(5R,8R,8aS)-8-Methyl-5-((Z)-non-...	263	C18H33N	1000367-14-0	47		
3	2,5-Cyclohexadiene-1,4-dione, 2-...	138	C7H6O3	000615-91-8	43		
4	Gephyrotoxin 235b'	235	C16H29N	120030-10-6	43		
5	Succinic acid, dodec-2-en-1-yl d...	422	C26H46O4	1010391-18-3	43		



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF053122\
 Data File : BF128602.D
 Acq On : 31 May 2022 18:56
 Operator : CG\JU
 Sample : N3109-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT 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
2-Furanmethanol...	5.610	4.6	ng	148328	1	6.845	641570	20.0
Dodecyl isobuty...	9.434	3.7	ng	143356	3	9.881	783488	20.0
unknown10.781	10.781	8.6	ng	309444	4	11.369	718878	20.0
unknown11.010	11.010	5.4	ng	194332	4	11.369	718878	20.0
unknown11.880	11.880	2.8	ng	101202	4	11.369	718878	20.0
n-Hexadecanoic ...	11.910	4.6	ng	163789	4	11.369	718878	20.0
unknown11.980	11.980	9.3	ng	333951	4	11.369	718878	20.0
unknown12.151	12.151	17.1	ng	615819	4	11.369	718878	20.0
Ethanol, 2-[2-(...	13.033	13.5	ng	442367	5	14.016	655714	20.0
unknown13.357	13.357	26.9	ng	882031	5	14.016	655714	20.0
unknown13.533	13.533	9.7	ng	316241	5	14.016	655714	20.0
2,3-Dimethylhyd...	13.586	42.5	ng	1395180	5	14.016	655714	20.0
Triphenylphosph...	14.121	10.7	ng	349323	5	14.016	655714	20.0
(5R,8R,8aS)-8-M...	14.286	42.3	ng	1386730	5	14.016	655714	20.0
unknown14.974	14.974	34.2	ng	893427	6	15.486	522037	20.0