

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122123\
 Data File : BF136670.D
 Acq On : 21 Dec 2023 15:43
 Operator : CG\JU
 Sample : 05832-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CHRT-20103

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-BF122023.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF136670.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.016	494	498	505	rVB	23833	26880	1.23%	0.205%
2	5.422	563	567	570	rBV	1324974	1256879	57.68%	9.568%
3	5.998	662	665	672	rBV	56381	67531	3.10%	0.514%
4	6.057	672	675	689	rVB	86036	106137	4.87%	0.808%
5	6.422	733	737	749	rBV	1253385	1251067	57.41%	9.524%
6	6.781	795	798	815	rBV2	582387	683015	31.34%	5.200%
7	6.957	820	828	831	rVV2	30437	65251	2.99%	0.497%
8	6.998	831	835	851	rVB	27906	106117	4.87%	0.808%
9	7.340	889	893	896	rBV	831965	761819	34.96%	5.800%
10	7.745	958	962	966	rBV4	18150	28521	1.31%	0.217%
11	7.887	982	986	996	rVB3	44563	78957	3.62%	0.601%
12	8.016	1003	1008	1011	rBV3	20607	29793	1.37%	0.227%
13	8.057	1011	1015	1018	rBV	614873	499830	22.94%	3.805%
14	8.416	1073	1076	1083	rVB4	32930	40813	1.87%	0.311%
15	8.639	1111	1114	1119	rBV3	33581	40076	1.84%	0.305%
16	9.051	1181	1184	1187	rBV3	25854	32688	1.50%	0.249%
17	9.133	1194	1198	1201	rBV	1505180	1251661	57.44%	9.529%
18	9.216	1207	1212	1215	rBV3	51932	70345	3.23%	0.536%
19	9.475	1253	1256	1258	rBV3	23525	28594	1.31%	0.218%
20	9.522	1261	1264	1265	rBV2	56387	50929	2.34%	0.388%
21	9.628	1280	1282	1286	rBV4	34491	31427	1.44%	0.239%
22	9.681	1287	1291	1293	rBV2	61469	75702	3.47%	0.576%
23	9.722	1295	1298	1301	rVB2	129973	113674	5.22%	0.865%
24	9.757	1301	1304	1307	rBV3	38247	62704	2.88%	0.477%
25	9.816	1307	1314	1317	rBV	526789	563971	25.88%	4.293%
26	9.928	1330	1333	1335	rVB4	43940	38935	1.79%	0.296%
27	10.075	1355	1358	1361	rBV2	69735	84444	3.88%	0.643%
28	10.222	1380	1383	1386	rBV2	132518	126515	5.81%	0.963%
29	10.604	1445	1448	1452	rBV	562410	541372	24.84%	4.121%
30	11.204	1548	1550	1558	rVB2	217569	278555	12.78%	2.121%
31	11.310	1565	1568	1570	rBV	406947	371233	17.04%	2.826%
32	11.845	1656	1659	1662	rBV	269863	276445	12.69%	2.104%
33	12.022	1687	1689	1691	rBV	139607	122741	5.63%	0.934%
34	12.304	1734	1737	1739	rBV	208434	223585	10.26%	1.702%
35	12.410	1753	1755	1759	rBV	211745	258103	11.84%	1.965%
36	12.557	1777	1780	1785	rBV	237872	422186	19.37%	3.214%

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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-BF122023.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	12.757	1812	1814	1817	rVB2	278012	221474	10.16%	1.686%
38	12.786	1817	1819	1821	rBV	282038	270691	12.42%	2.061%
39	12.910	1836	1840	1849	rBV2	1335983	2179116	100.00%	16.589%
40	13.274	1900	1902	1906	rBV	370265	396133	18.18%	3.016%

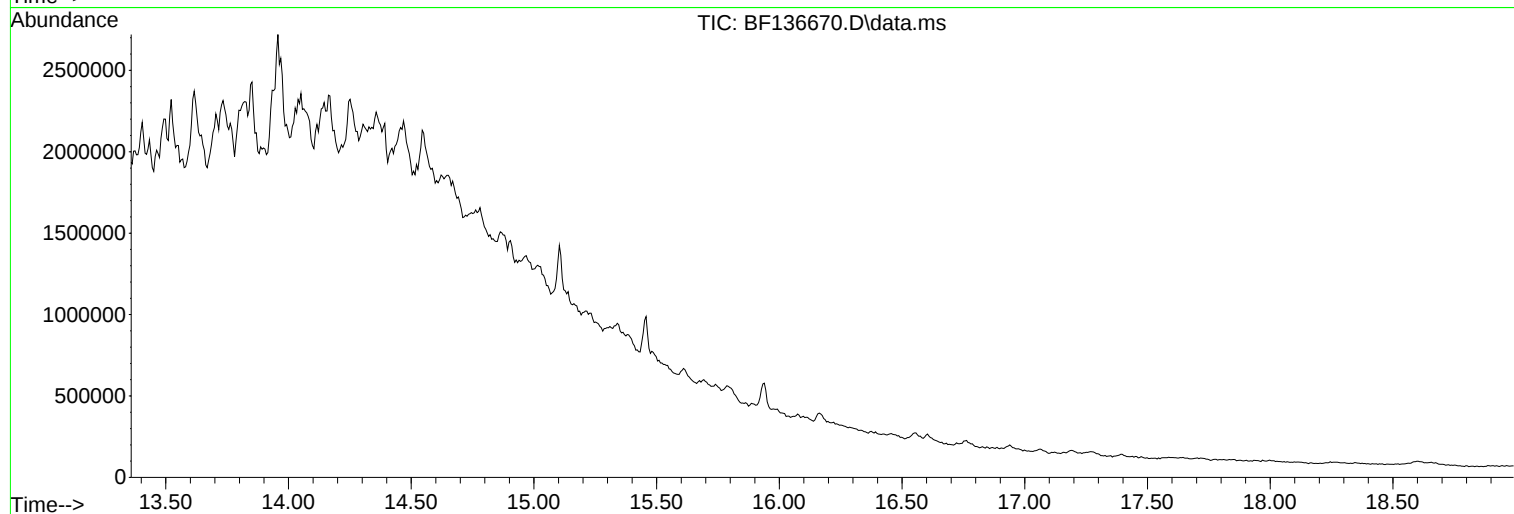
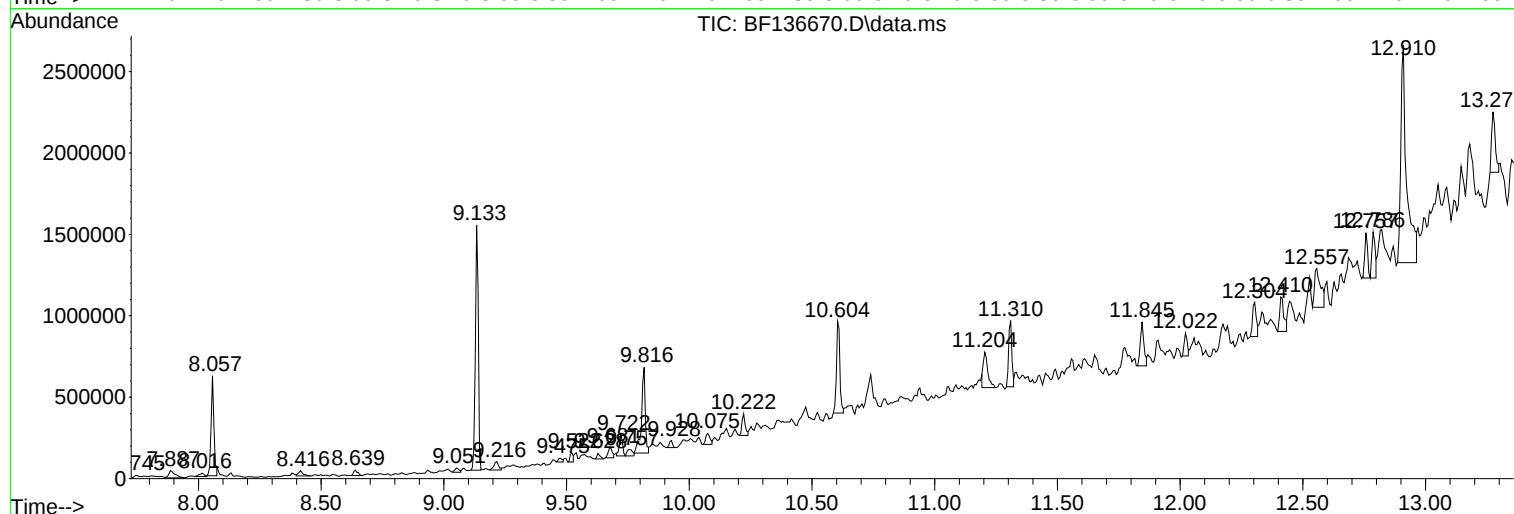
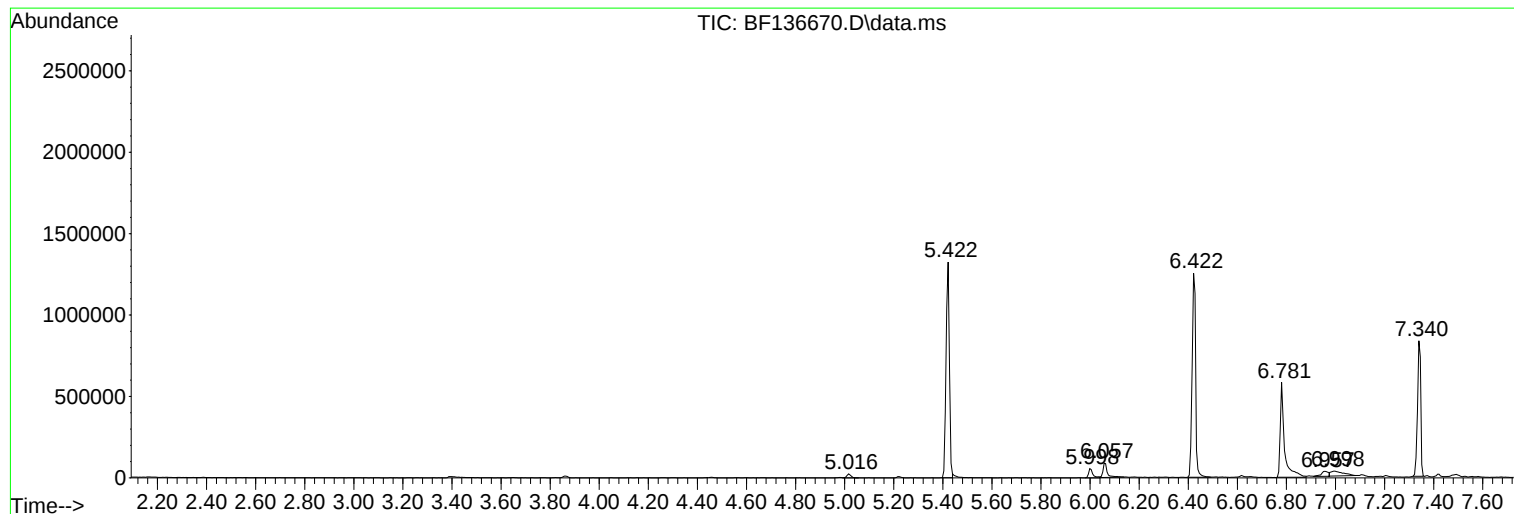
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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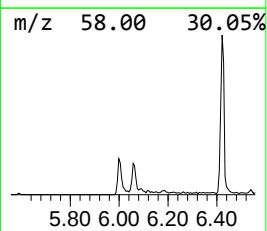
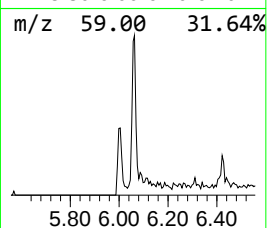
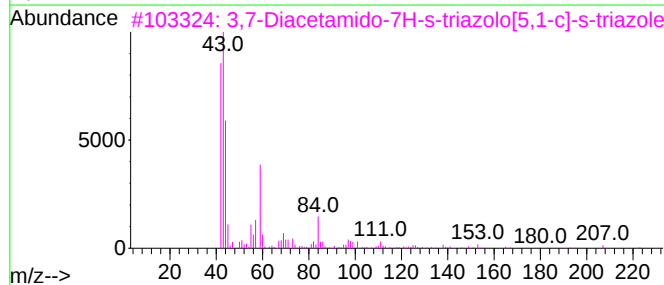
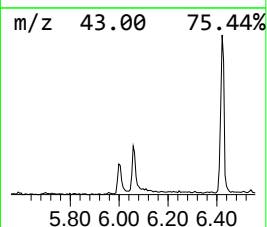
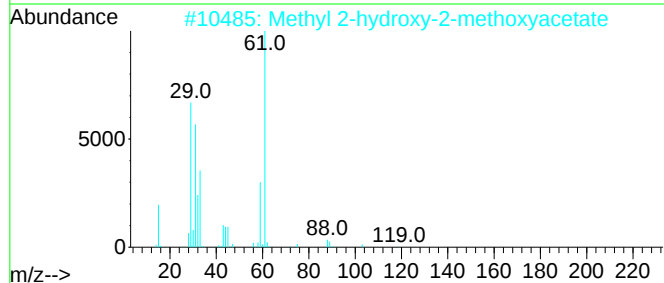
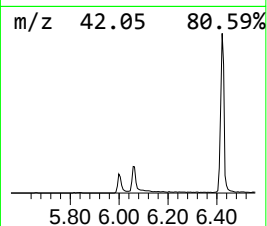
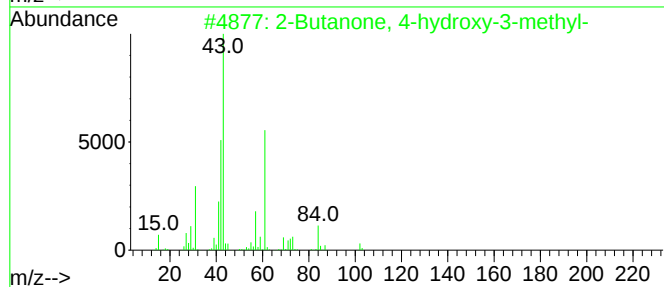
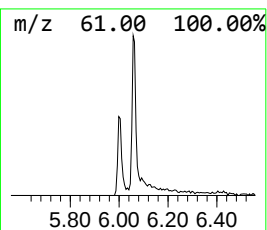
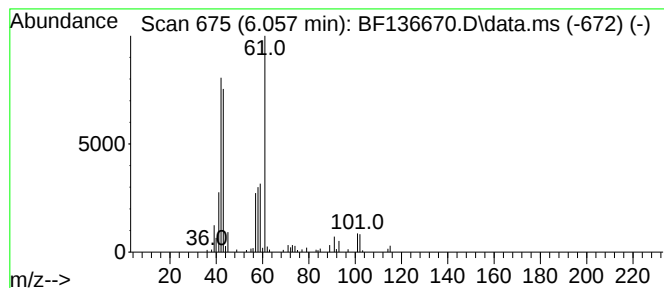
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.057 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.057	3.11 ng	106137	1,4-Dichlorobenzene-d4	6.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanone, 4-hydroxy-3-methyl-	102	C5H10O2	003393-64-4	42		
2	Methyl 2-hydroxy-2-methoxyacetate	120	C4H8O4	019757-97-2	25		
3	3,7-Diacetamido-7H-s-triazolo[5,...	223	C7H9N7O2	1000211-49-1	23		
4	N-Methoxy-N-methylacetamide	103	C4H9NO2	078191-00-1	17		
5	2-Nitro-tertiary butanol	119	C4H9NO3	005447-98-3	13		



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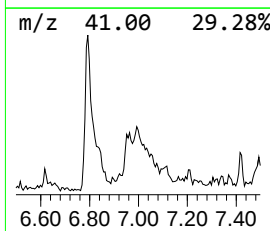
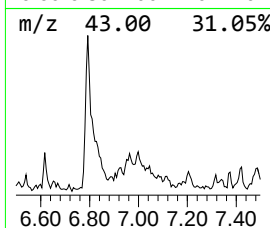
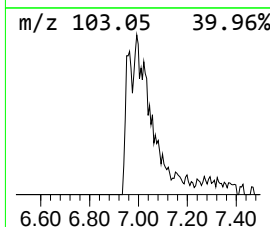
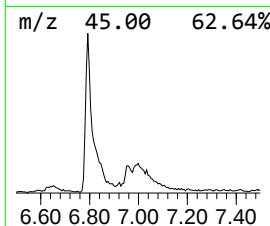
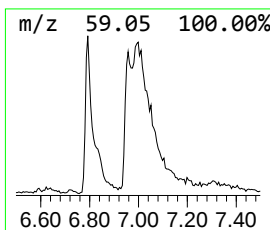
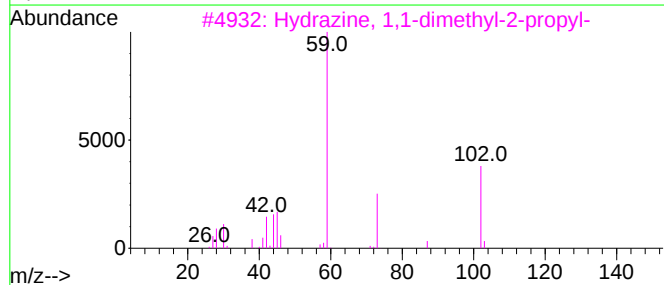
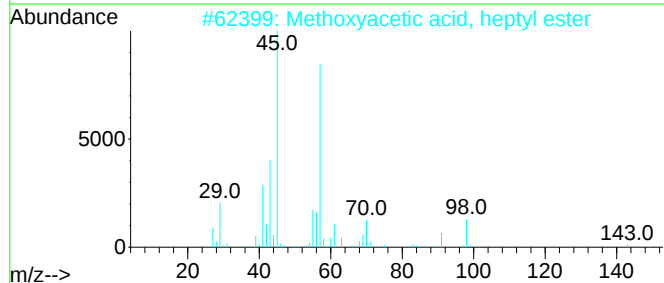
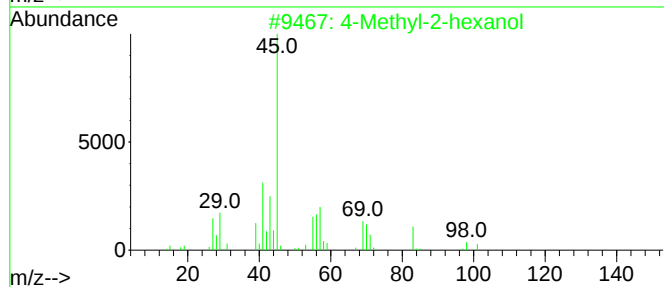
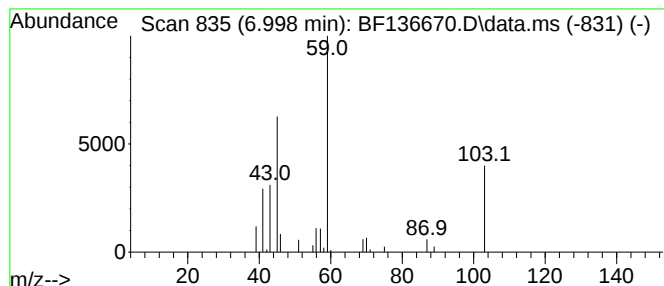
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.998 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.998	3.11 ng	106117	1,4-Dichlorobenzene-d4	6.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Methyl-2-hexanol	116	C7H16O	002313-61-3	9
2			Methoxyacetic acid, heptyl ester	188	C10H20O3	168920-36-3	9
3			Hydrazine, 1,1-dimethyl-2-propyl-	102	C5H14N2	052728-54-8	9
4			Propanoic acid, 2-nitro-, methyl...	133	C4H7NO4	006118-50-9	9
5			Silane, dimethyl-	60	C2H8Si	001111-74-6	9



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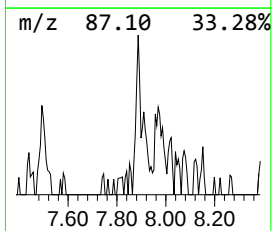
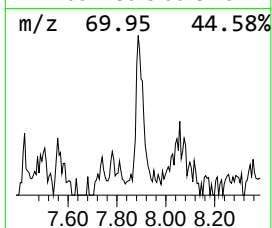
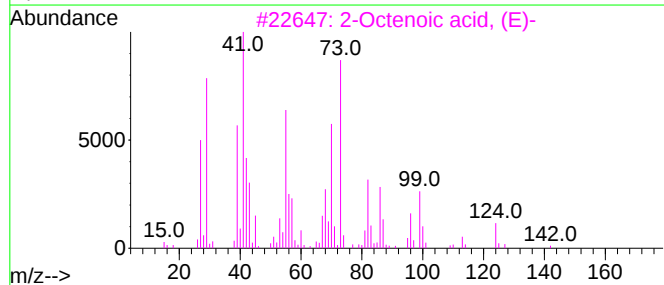
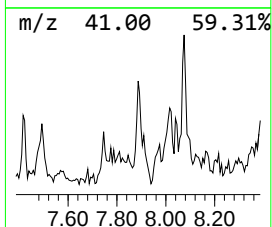
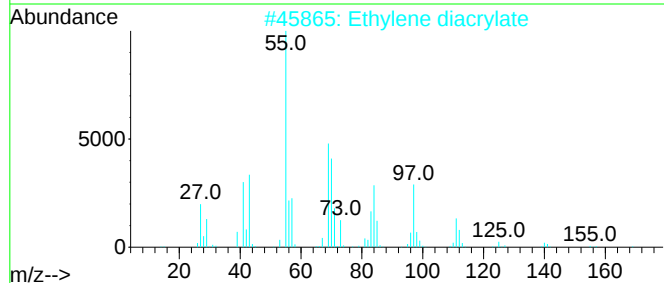
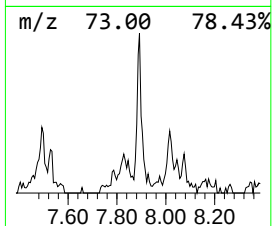
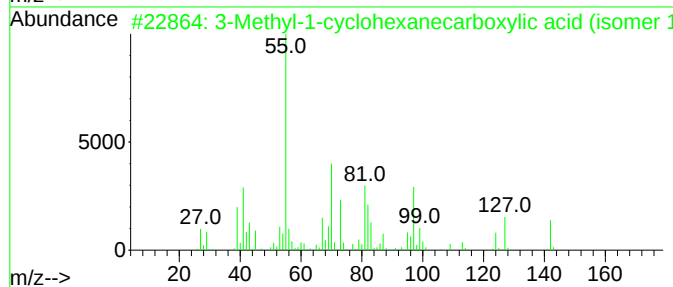
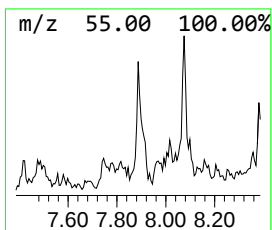
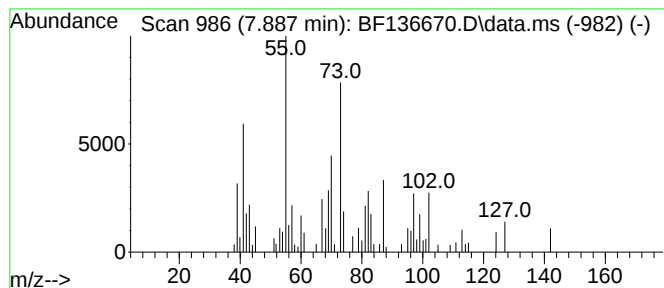
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 3-Methyl-1-cyclohexanecarbo... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.887	3.16 ng	78957	Naphthalene-d8	8.057

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Methyl-1-cyclohexanecarboxylic...	142	C8H14O2	1000453-57-0	64
2			Ethylene diacrylate	170	C8H10O4	002274-11-5	27
3			2-Octenoic acid, (E)-	142	C8H14O2	001871-67-6	16
4			2-Decene, (E)-	140	C10H20	020063-97-2	14
5			Cyclotetradecane	196	C14H28	000295-17-0	12



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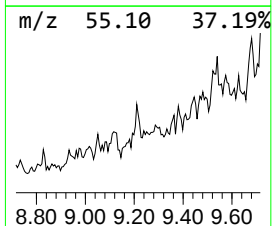
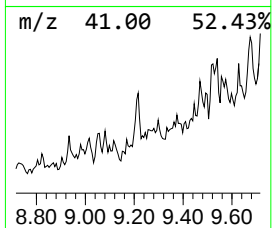
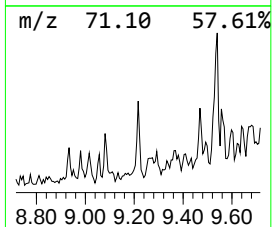
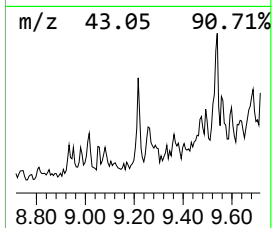
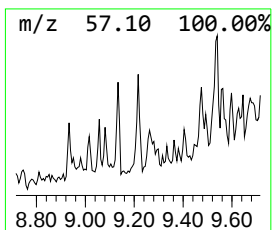
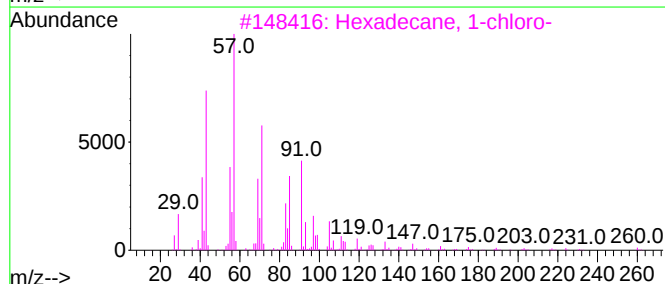
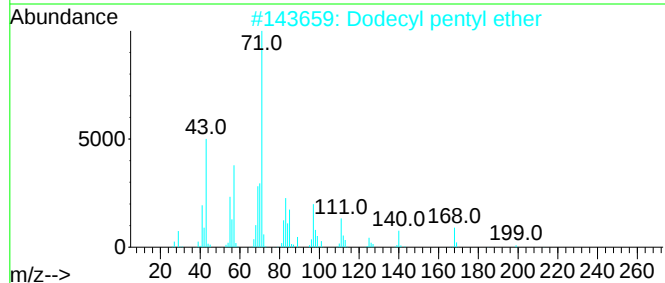
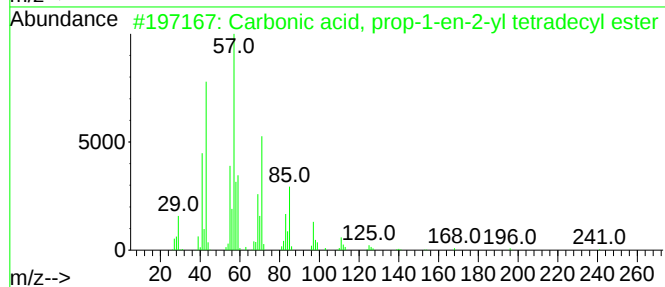
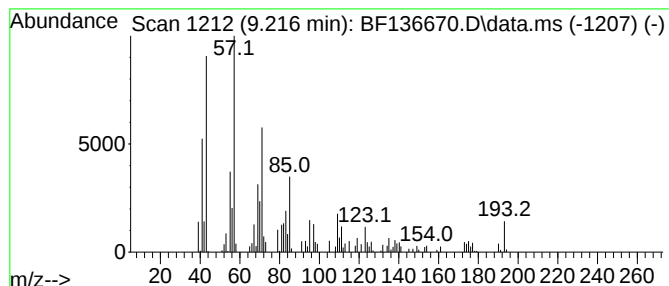
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 Carbonic acid, prop-1-en-2-... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.216	2.49 ng	70345	Acenaphthene-d10	9.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, prop-1-en-2-yl te...	298	C18H34O3	1000382-90-9	58
2			Dodecyl pentyl ether	256	C17H36O	1010406-41-6	50
3			Hexadecane, 1-chloro-	260	C16H33Cl	004860-03-1	50
4			2-Methyltetracosane	352	C25H52	001560-78-7	50
5			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	49



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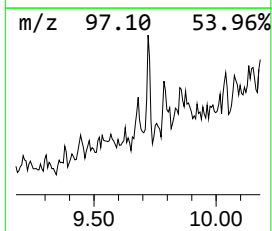
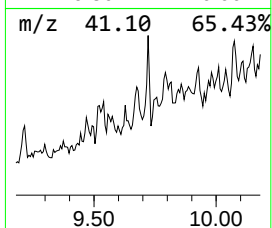
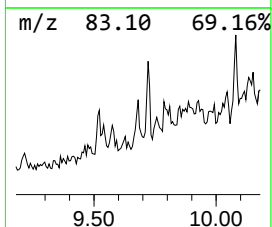
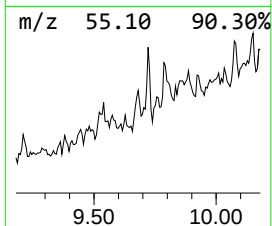
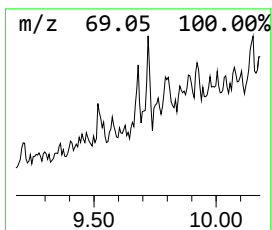
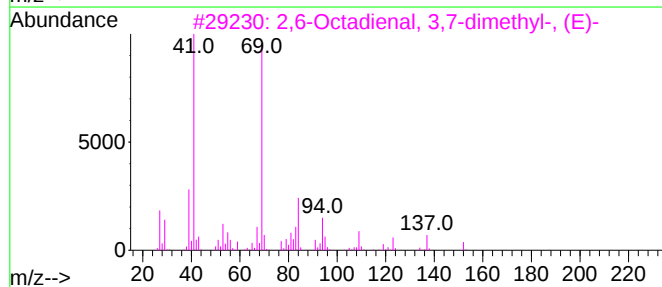
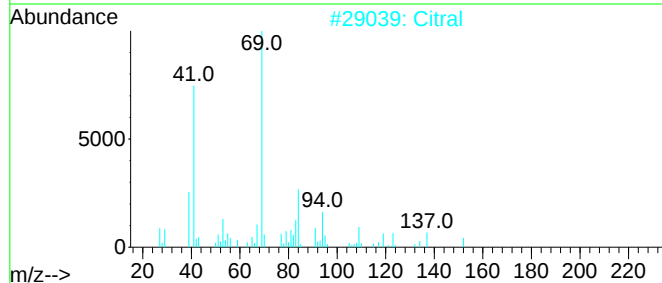
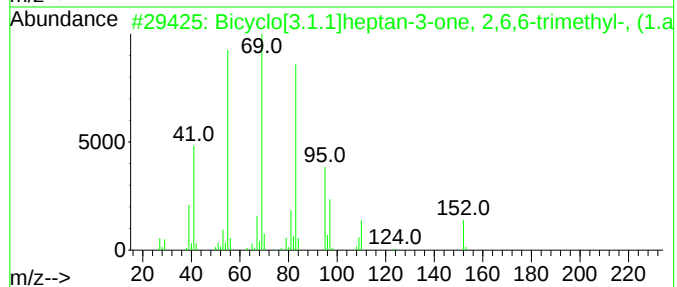
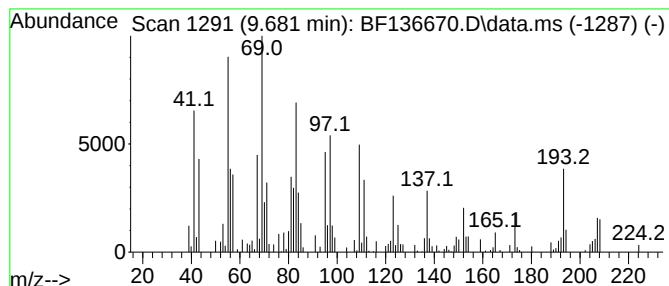
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 unknown9.681 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.681	2.68 ng	75702	Acenaphthene-d10	9.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[3.1.1]heptan-3-one, 2,6,...	152	C10H16O	015358-88-0	43
2			Citral	152	C10H16O	005392-40-5	35
3			2,6-Octadienal, 3,7-dimethyl-, (E)-	152	C10H16O	000141-27-5	35
4			1,1'-Bicyclohexyl, 2-ethyl-, cis-	194	C14H26	050991-12-3	30
5			Cyclohexane, 1-(cyclohexylmethyl)...	208	C15H28	054934-94-0	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122123\
Data File : BF136670.D
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Sample : 05832-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

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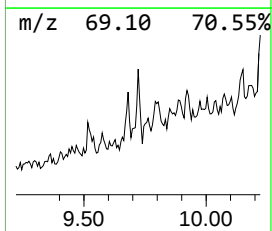
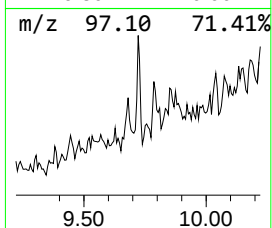
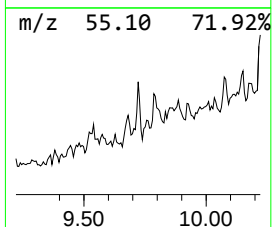
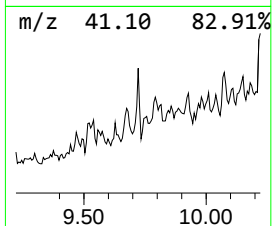
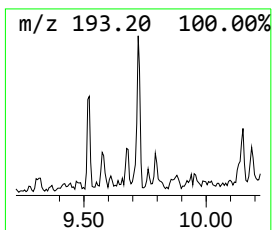
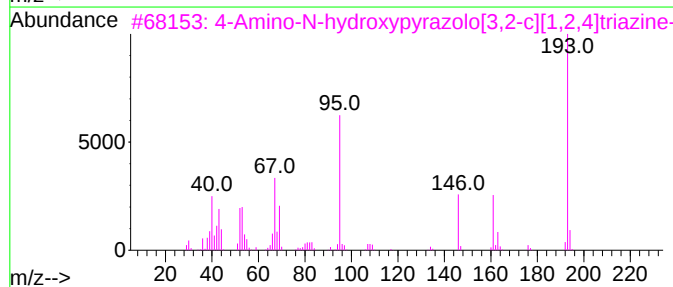
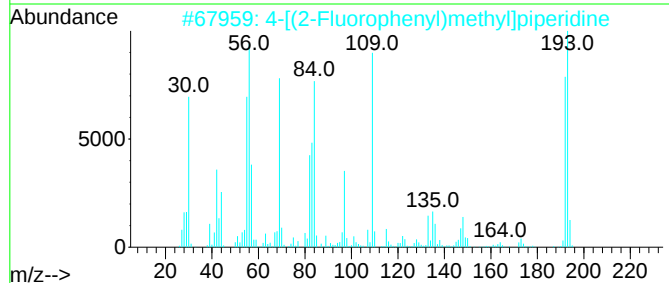
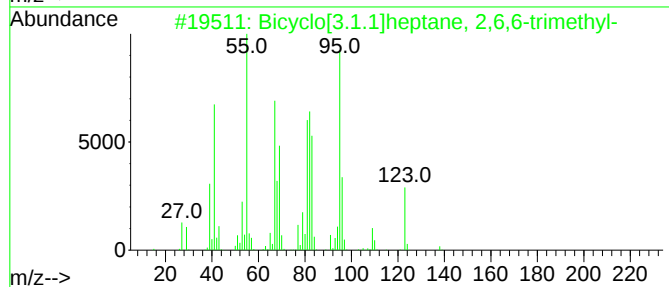
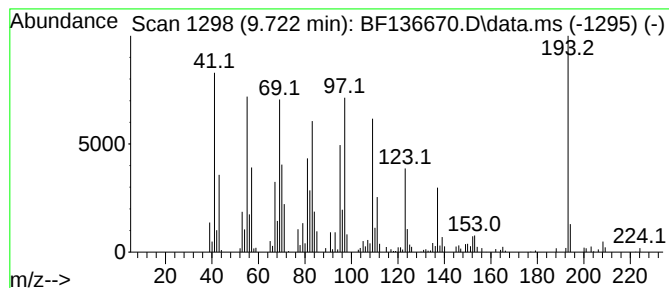
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown9.722 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.722	4.03 ng	113674	Acenaphthene-d10	9.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[3.1.1]heptane, 2,6,6-tri...	138	C10H18	000473-55-2	38
2			4-[(2-Fluorophenyl)methyl]piperi...	193	C12H16FN	194288-97-6	35
3			4-Amino-N-hydroxypyrazolo[3,2-c]...	193	C6H7N7O	1000443-50-2	25
4			Cyclohexane, 1-(cyclohexylmethyl...	194	C14H26	054823-96-0	18
5			N-(4-Fluorophenyl)piperidin-4-amine	194	C11H15FN2	038043-08-2	15



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122123\
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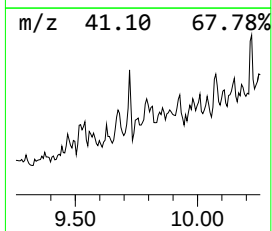
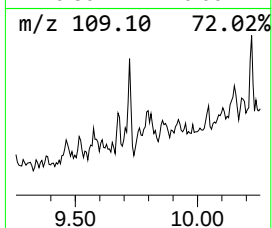
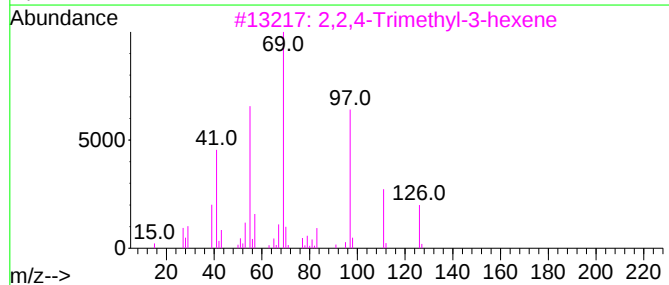
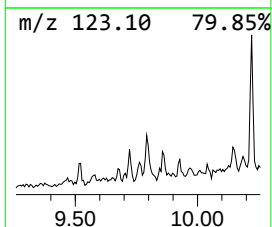
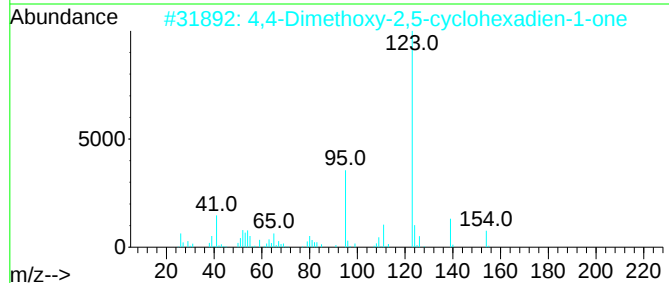
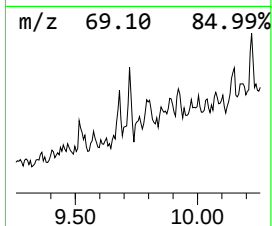
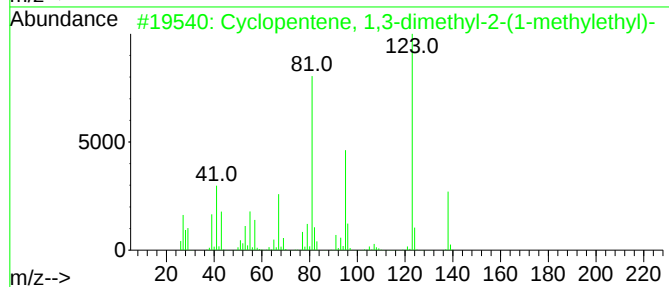
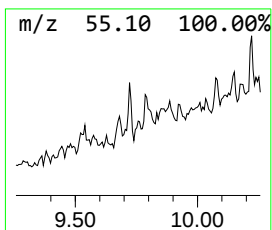
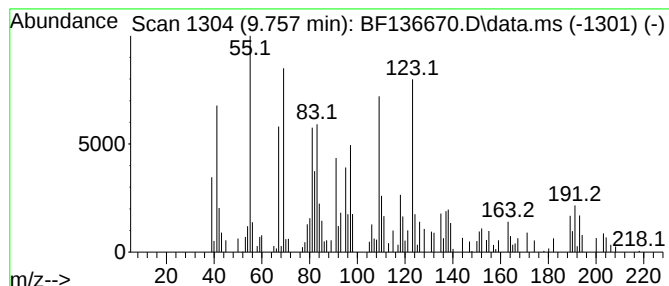
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown9.757 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.757	2.22 ng	62704	Acenaphthene-d10	9.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentene, 1,3-dimethyl-2-(1-...	138	C10H18	061142-32-3	25
2			4,4-Dimethoxy-2,5-cyclohexadien-...	154	C8H10O3	000935-50-2	18
3			2,2,4-Trimethyl-3-hexene	126	C9H18	059643-72-0	18
4			Cyclopropanemethanol, 2,2-dimeth...	154	C10H18O	005617-92-5	15
5			3-Fluoropropiophenone	152	C9H9FO	000455-67-4	14



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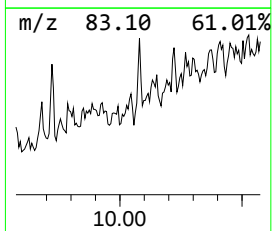
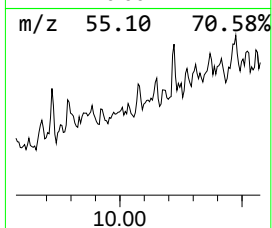
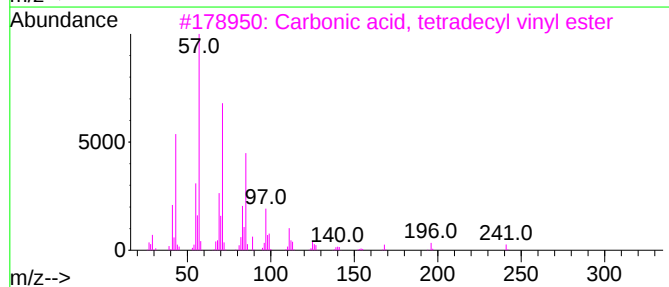
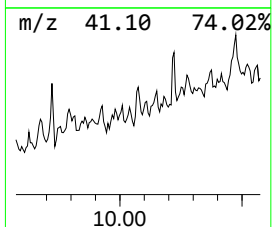
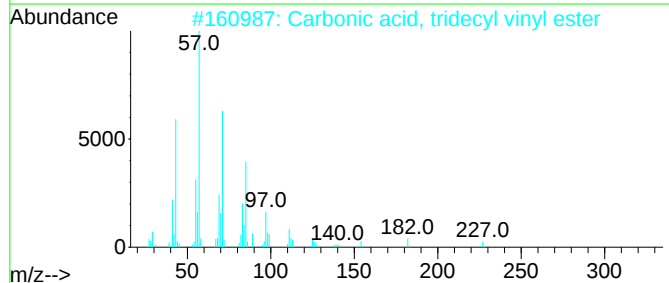
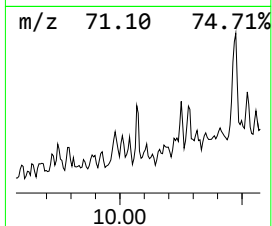
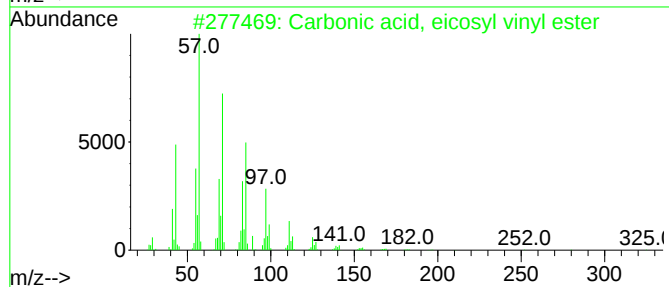
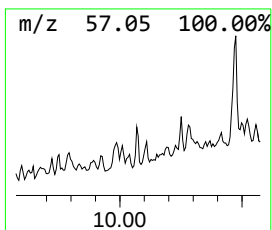
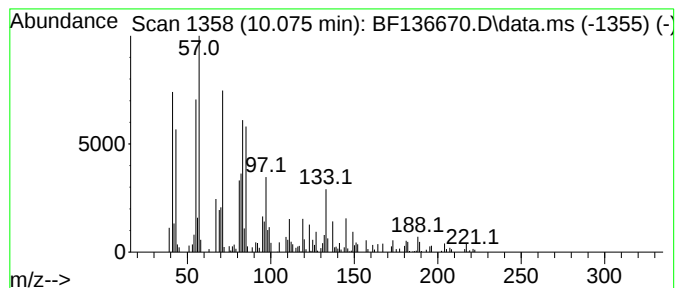
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Carbonic acid, eicosyl vinyl... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.075	2.99 ng	84444	Acenaphthene-d10	9.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	58
2			Carbonic acid, tridecyl vinyl ester	270	C16H30O3	1000382-54-7	49
3			Carbonic acid, tetradecyl vinyl ...	284	C17H32O3	1000382-54-5	46
4			Oxalic acid, cyclobutyl heptadec...	382	C23H42O4	1000309-70-7	46
5			Decane, 2,3,6-trimethyl-	184	C13H28	062238-12-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122123\
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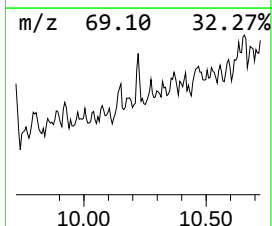
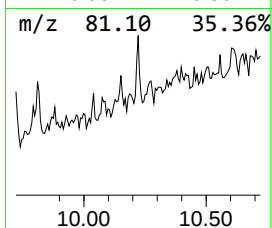
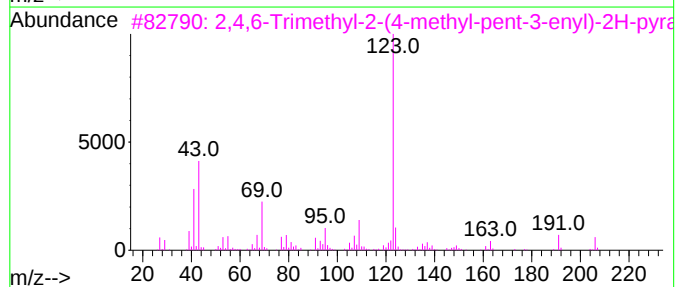
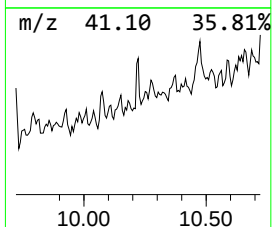
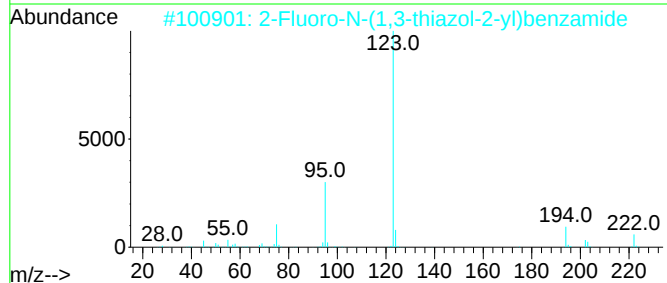
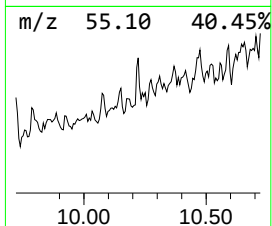
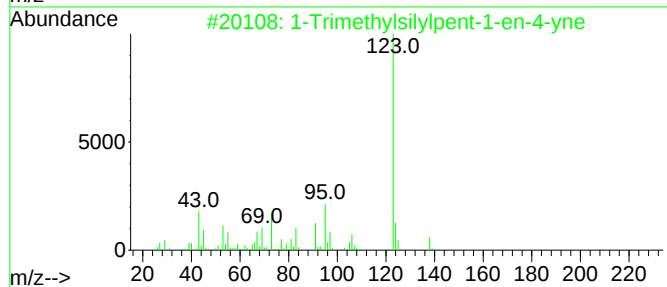
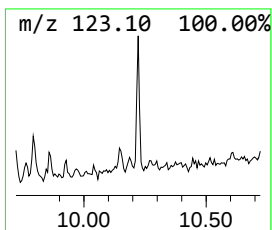
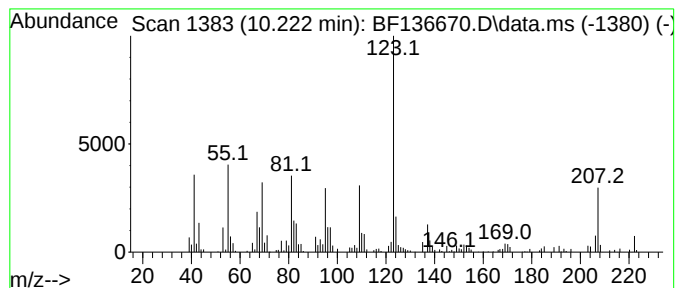
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 1-Trimethylsilylpent-1-en-4... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.222	4.49 ng	126515	Acenaphthene-d10	9.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Trimethylsilylpent-1-en-4-yne	138	C8H14Si	079516-25-9	50
2			2-Fluoro-N-(1,3-thiazol-2-yl)ben...	222	C10H7FN2OS	000319-38-0	49
3			2,4,6-Trimethyl-2-(4-methyl-pent...	206	C14H22O	1000193-58-6	49
4			2-Amino-4-acetamino anisole	180	C9H12N2O2	006375-47-9	47
5			Benzamide, 3-fluoro-N-butyl-N-et...	223	C13H18FNO	1010415-85-9	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122123\
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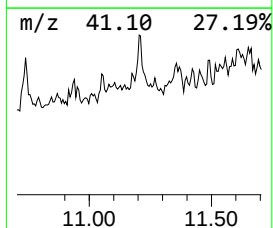
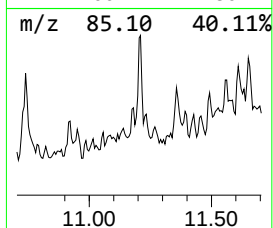
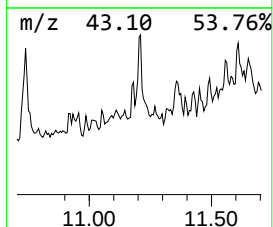
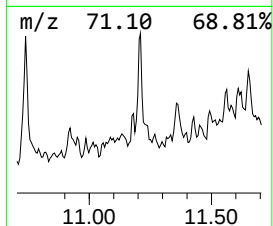
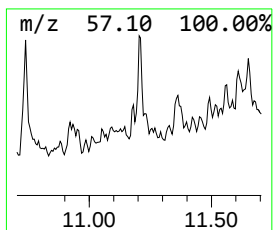
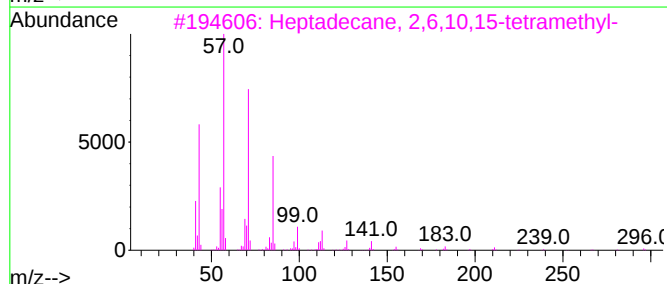
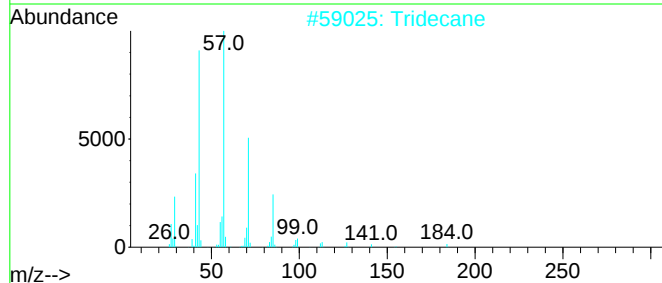
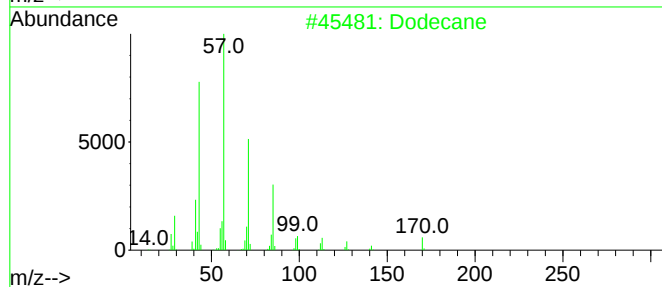
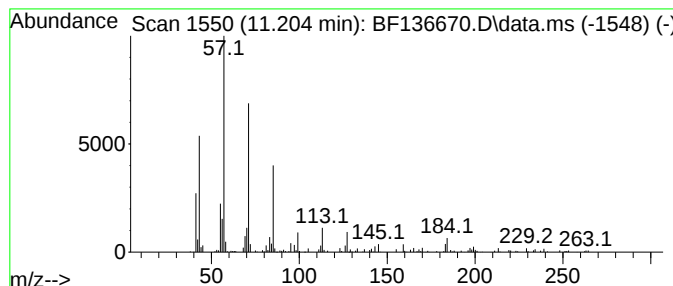
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Dodecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.204	15.01 ng	278555	Phenanthrene-d10	11.310

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Dodecane	170	C12H26	000112-40-3	90
2		Tridecane	184	C13H28	000629-50-5	87
3		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	80
4		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	72
5		Decane, 3,6-dimethyl-	170	C12H26	017312-53-7	72



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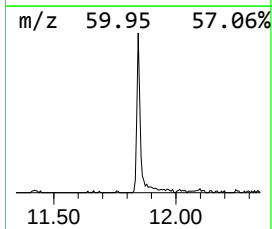
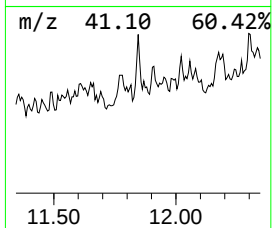
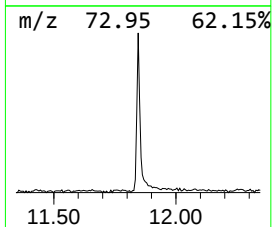
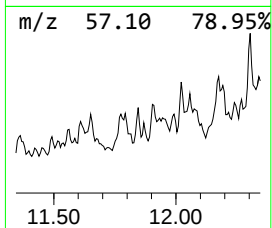
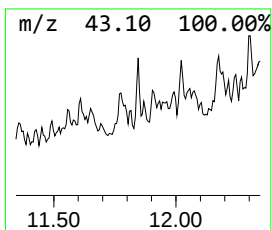
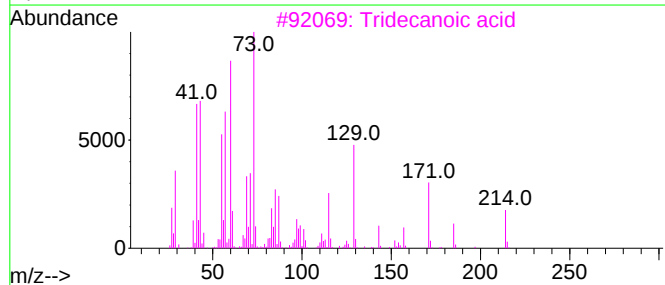
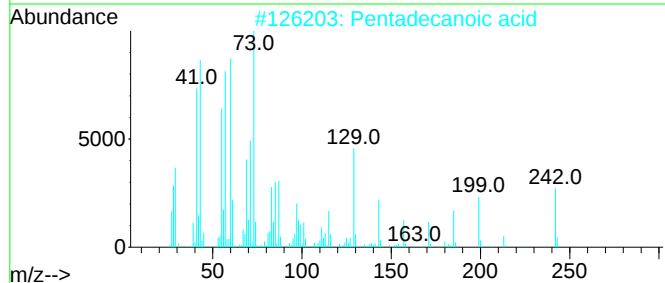
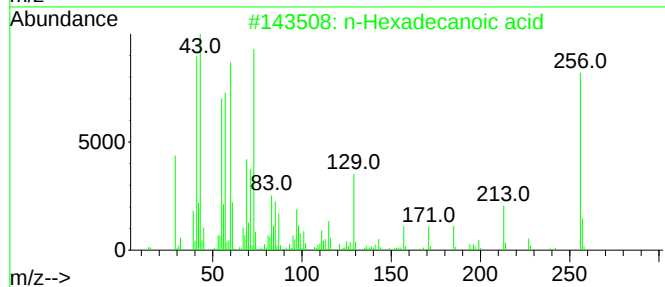
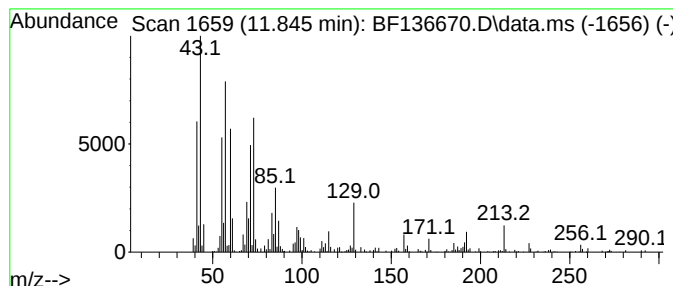
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.845	14.89 ng	276445	Phenanthrene-d10	11.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	70
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	62
3			Tridecanoic acid	214	C13H26O2	000638-53-9	60
4			Sulfurous acid, 2-propyl tetra...	320	C17H36O3S	1010309-12-5	41
5			Nonyl tetradecyl ether	340	C23H48O	1000406-37-6	38



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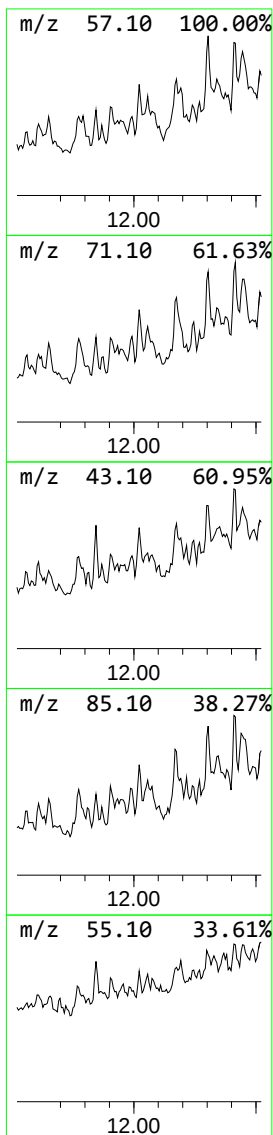
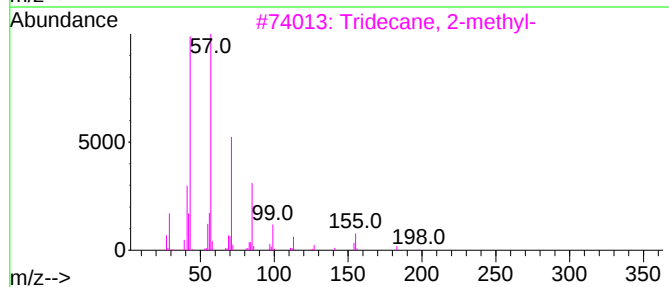
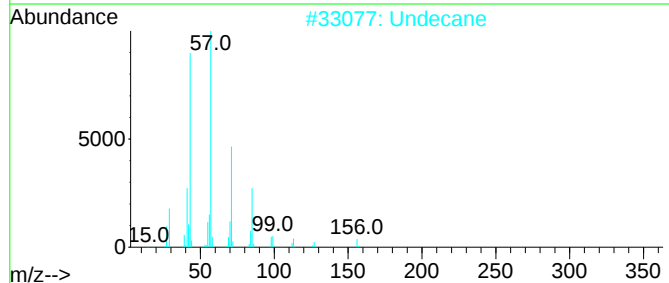
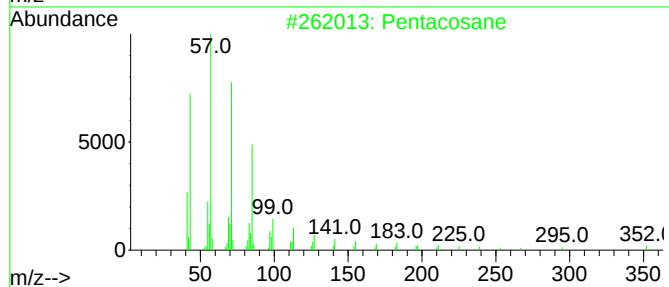
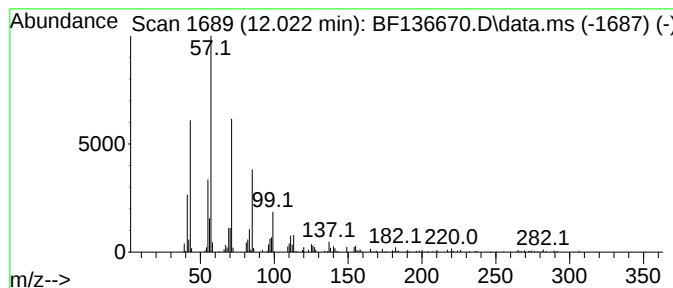
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Pentacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.022	6.61 ng	122741	Phenanthrene-d10	11.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentacosane	352	C25H52	000629-99-2	83
2			Undecane	156	C11H24	001120-21-4	81
3			Tridecane, 2-methyl-	198	C14H30	001560-96-9	80
4			Heptadecane	240	C17H36	000629-78-7	80
5			Pentadecane, 6-methyl-	226	C16H34	010105-38-1	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122123\
Data File : BF136670.D
Acq On : 21 Dec 2023 15:43
Operator : CG\JU
Sample : 05832-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CHRT-20103

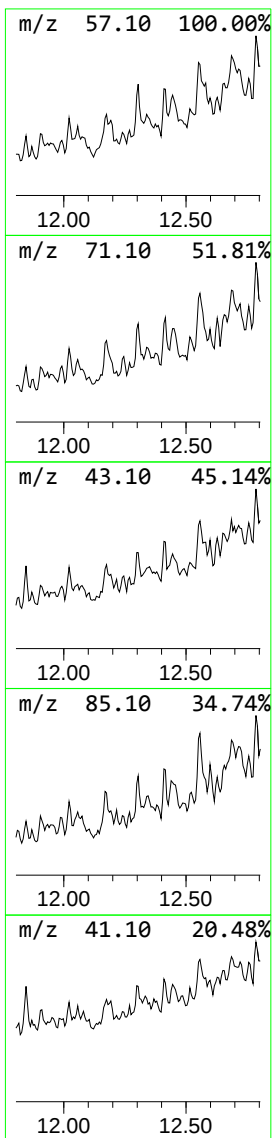
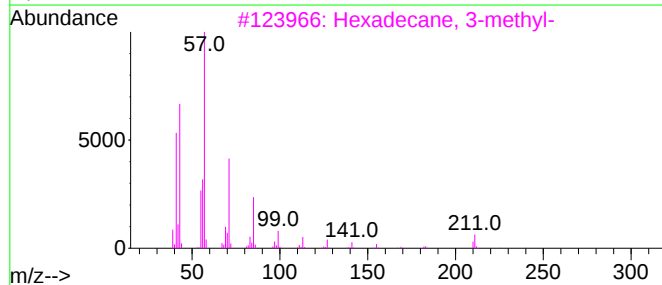
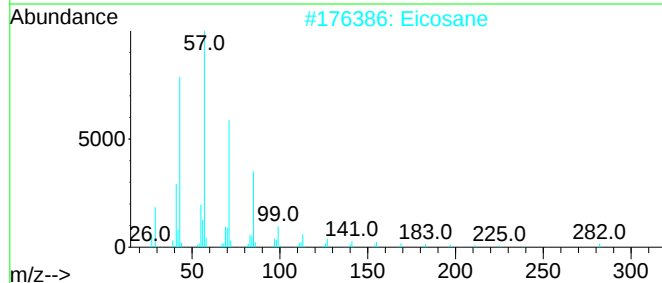
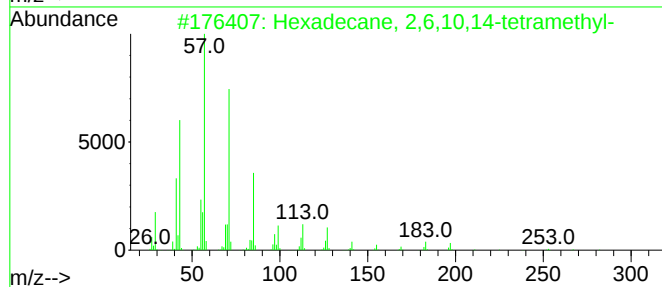
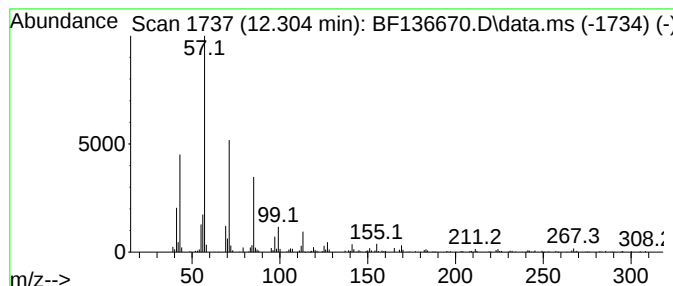
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Hexadecane, 2,6,10,14-tetra... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.304	12.05 ng	223585	Phenanthrene-d10	11.310

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	90
2		Eicosane	282	C20H42	000112-95-8	90
3		Hexadecane, 3-methyl-	240	C17H36	006418-43-5	90
4		Nonadecane, 2-methyl-	282	C20H42	001560-86-7	90
5		Tetradecane, 3-methyl-	212	C15H32	018435-22-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122123\
Data File : BF136670.D
Acq On : 21 Dec 2023 15:43
Operator : CG\JU
Sample : 05832-01
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Instrument :
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ClientSampleId :
CHRT-20103

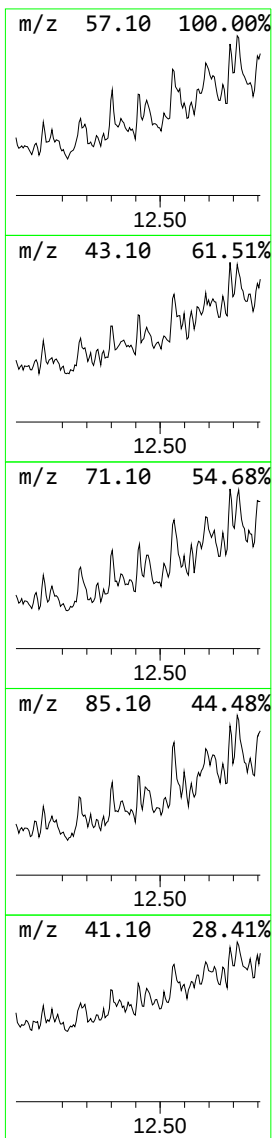
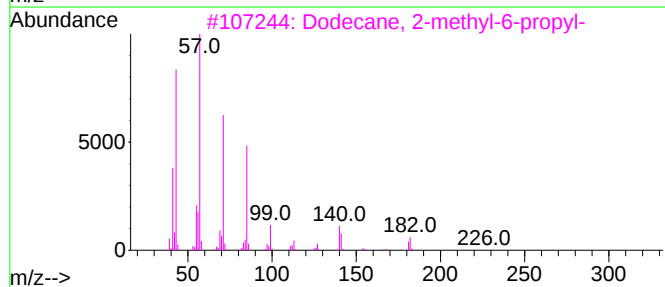
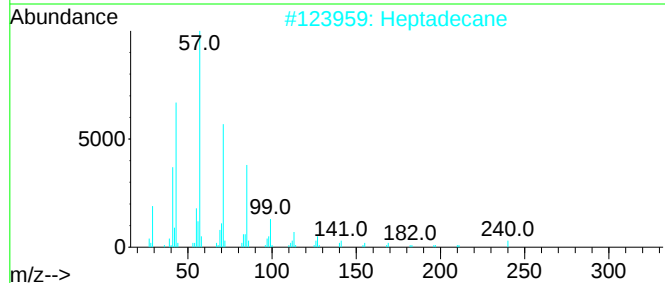
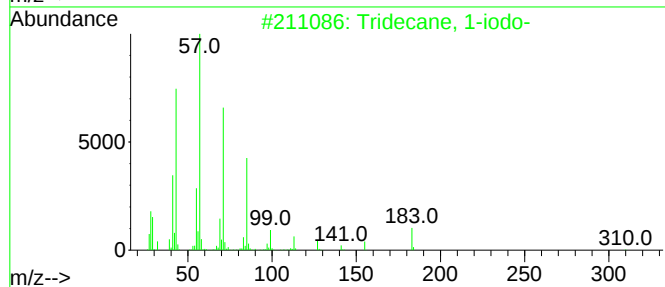
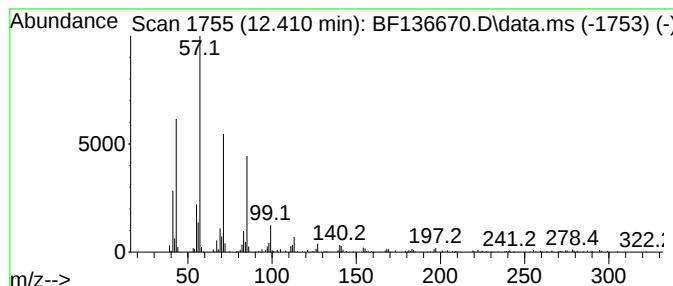
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Tridecane, 1-iodo- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.410	13.91 ng	258103	Phenanthrene-d10	11.310

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	90
2		Heptadecane	240	C17H36	000629-78-7	87
3		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	86
4		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	68
5		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122123\
Data File : BF136670.D
Acq On : 21 Dec 2023 15:43
Operator : CG\JU
Sample : 05832-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

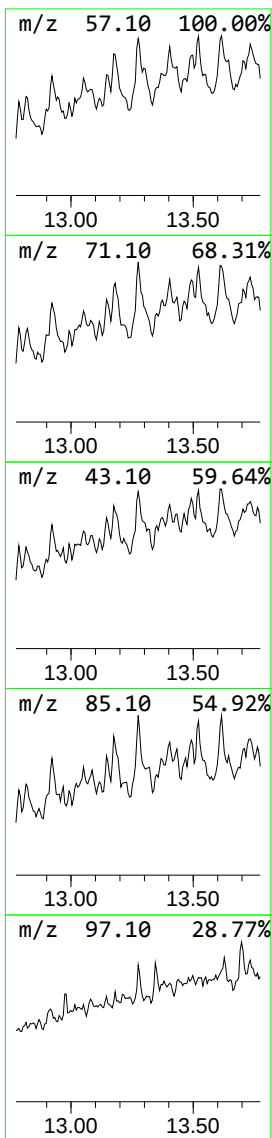
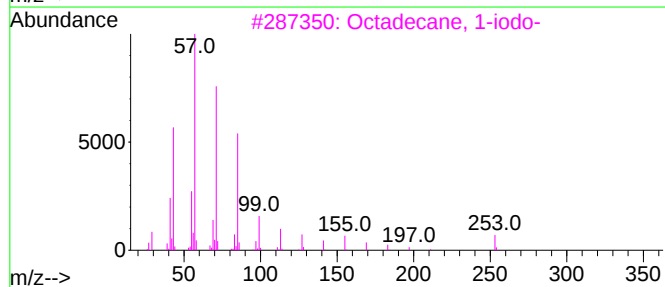
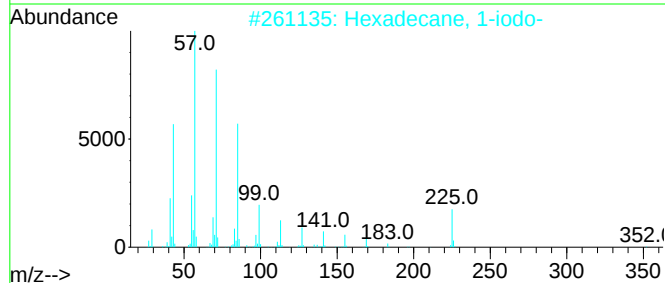
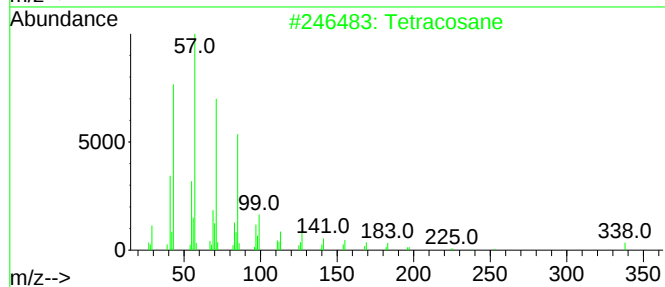
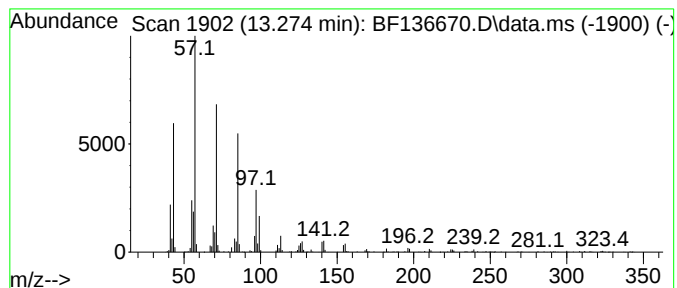
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Tetracosane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.274	20.00 ng	396133	Chrysene-d12	13.969	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	86
2	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	80
3	Octadecane, 1-iodo-	380	C18H37I	000629-93-6	72
4	Octadecane, 4-methyl-	268	C19H40	010544-95-3	58
5	Octadecane, 2-methyl-	268	C19H40	001560-88-9	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122123\
Data File : BF136670.D
Acq On : 21 Dec 2023 15:43
Operator : CG\JU
Sample : 05832-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CHRT-20103

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.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--			
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unknown6.998	6.998	3.1	ng	106117	1	6.781	683015	20.0
3-Methyl-1-cycl...	7.887	3.2	ng	78957	2	8.057	499830	20.0
Carbonic acid, ...	9.216	2.5	ng	70345	3	9.816	563971	20.0
unknown9.681	9.681	2.7	ng	75702	3	9.816	563971	20.0
unknown9.722	9.722	4.0	ng	113674	3	9.816	563971	20.0
unknown9.757	9.757	2.2	ng	62704	3	9.816	563971	20.0
Carbonic acid, ...	10.075	3.0	ng	84444	3	9.816	563971	20.0
1-Trimethylsily...	10.222	4.5	ng	126515	3	9.816	563971	20.0
Dodecane	11.204	15.0	ng	278555	4	11.310	371233	20.0
n-Hexadecanoic ...	11.845	14.9	ng	276445	4	11.310	371233	20.0
Pentacosane	12.022	6.6	ng	122741	4	11.310	371233	20.0
Hexadecane, 2,6...	12.304	12.1	ng	223585	4	11.310	371233	20.0
Tridecane, 1-iodo-	12.410	13.9	ng	258103	4	11.310	371233	20.0
Tetracosane	13.274	20.0	ng	396133	5	13.969	396133	20.0