

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050322\
 Data File : BF128047.D
 Acq On : 03 May 2022 15:52
 Operator : CG\JU
 Sample : N2650-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ETGI-362

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-BF042222.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF128047.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.475	197	202	209	rBV	59491	95291	3.28%	0.288%
2	4.034	292	297	303	rVB	44601	54548	1.88%	0.165%
3	5.540	549	553	556	rVB	2632989	2446252	84.26%	7.393%
4	6.546	718	724	727	rBV	2453970	2704249	93.15%	8.173%
5	6.916	783	787	790	rBV	1554550	1307438	45.04%	3.951%
6	7.481	877	883	886	rBV	1931271	1847086	63.62%	5.582%
7	7.698	914	920	922	rBV2	52150	61489	2.12%	0.186%
8	8.092	984	987	992	rBV	91261	124762	4.30%	0.377%
9	8.198	1001	1005	1009	rVB	1942820	1746812	60.17%	5.279%
10	8.245	1010	1013	1016	rVB3	72786	63937	2.20%	0.193%
11	8.275	1016	1018	1021	rBV3	45282	47516	1.64%	0.144%
12	8.434	1041	1045	1047	rBV2	83875	92878	3.20%	0.281%
13	8.492	1051	1055	1058	rVB3	87935	113915	3.92%	0.344%
14	8.581	1066	1070	1072	rBV2	71208	87342	3.01%	0.264%
15	8.634	1077	1079	1081	rBV2	73696	80058	2.76%	0.242%
16	8.657	1081	1083	1086	rVB	148015	135202	4.66%	0.409%
17	8.698	1087	1090	1092	rBV2	86527	74153	2.55%	0.224%
18	8.728	1092	1095	1096	rBV2	66744	64597	2.23%	0.195%
19	8.781	1102	1104	1107	rVB3	76922	66730	2.30%	0.202%
20	8.816	1107	1110	1111	rBV2	112524	114983	3.96%	0.348%
21	8.839	1111	1114	1118	rVV3	115474	143045	4.93%	0.432%
22	8.887	1119	1122	1126	rVB4	144984	162709	5.60%	0.492%
23	8.939	1129	1131	1133	rVB2	134716	98234	3.38%	0.297%
24	8.969	1133	1136	1139	rBV4	87711	120830	4.16%	0.365%
25	9.022	1143	1145	1147	rBV2	150706	139772	4.81%	0.422%
26	9.075	1152	1154	1158	rBV5	118609	124454	4.29%	0.376%
27	9.163	1161	1169	1171	rBV7	272637	597915	20.60%	1.807%
28	9.192	1171	1174	1175	rBV2	259217	207150	7.14%	0.626%
29	9.275	1183	1188	1192	rBV	2989721	2903158	100.00%	8.774%
30	9.351	1197	1201	1204	rBV2	760744	855165	29.46%	2.585%
31	9.592	1239	1242	1243	rBV2	258699	286168	9.86%	0.865%
32	9.663	1251	1254	1258	rBV2	1209904	1175987	40.51%	3.554%
33	9.722	1262	1264	1268	rVB3	380676	409356	14.10%	1.237%
34	9.822	1278	1281	1283	rBV2	871015	873012	30.07%	2.638%
35	9.869	1286	1289	1292	rVB	1406798	1371958	47.26%	4.146%
36	9.910	1293	1296	1298	rBV3	382619	504463	17.38%	1.525%

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37	9.945	1299	1302	1303	rVV2	650170	652405	22.47%	1.972%
38	9.963	1303	1305	1308	rVB	1695441	1557165	53.64%	4.706%
39	9.992	1308	1310	1312	rBV3	298900	308673	10.63%	0.933%
40	10.281	1356	1359	1360	rBV2	380752	405493	13.97%	1.226%
41	10.369	1372	1374	1377	rVB	862278	705973	24.32%	2.134%
42	10.763	1437	1441	1444	rBV	1635576	1852272	63.80%	5.598%
43	10.875	1458	1460	1462	rVB	701499	538127	18.54%	1.626%
44	11.463	1557	1560	1563	rBV	1659150	1505997	51.87%	4.552%
45	13.045	1825	1829	1832	rVB	2081507	1805403	62.19%	5.456%
46	14.104	2005	2009	2012	rVB	1098071	1012973	34.89%	3.061%
47	14.574	2086	2089	2092	rVB	352604	305903	10.54%	0.925%
48	15.610	2260	2265	2272	rVB	887341	1134698	39.08%	3.429%

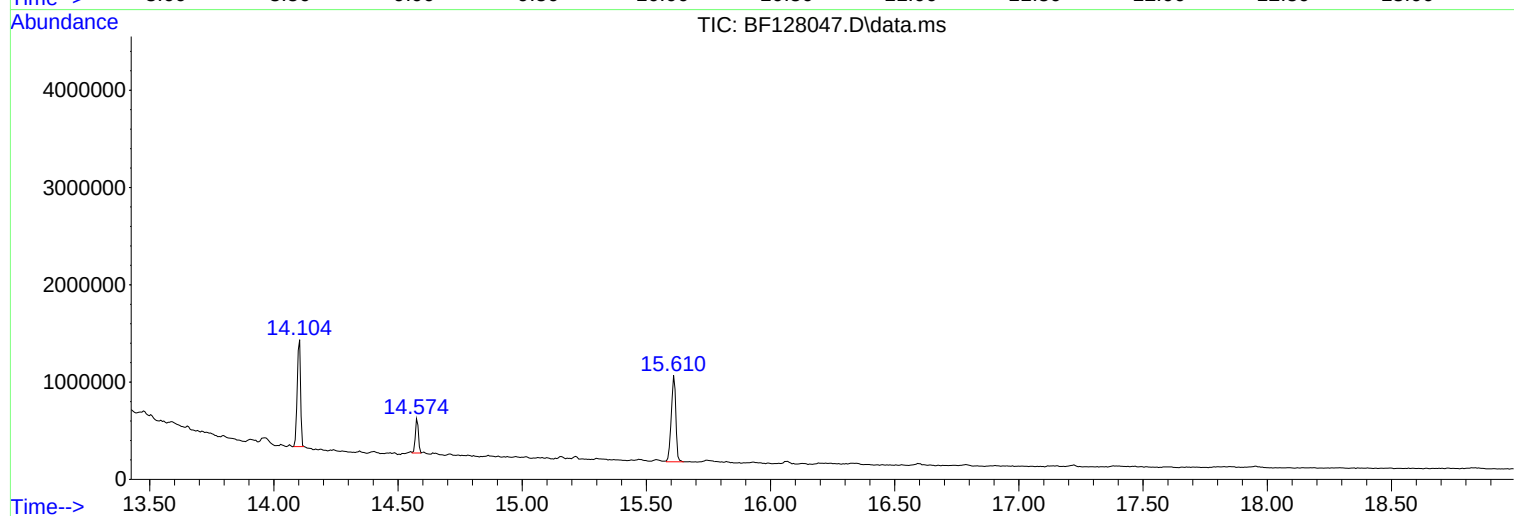
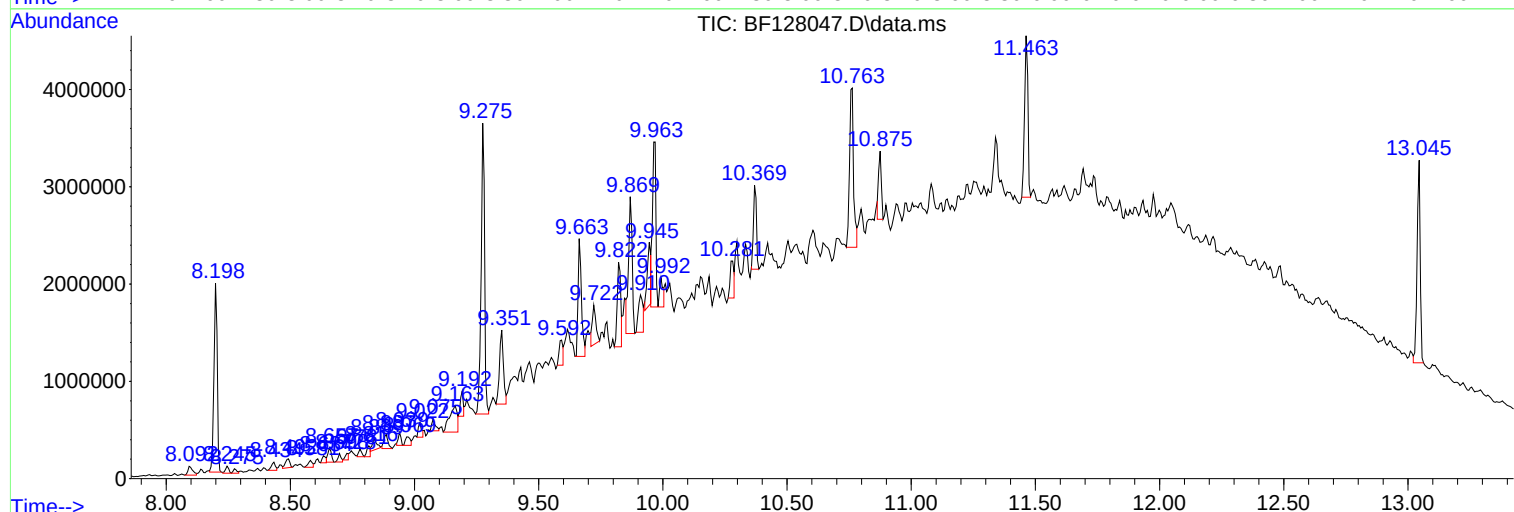
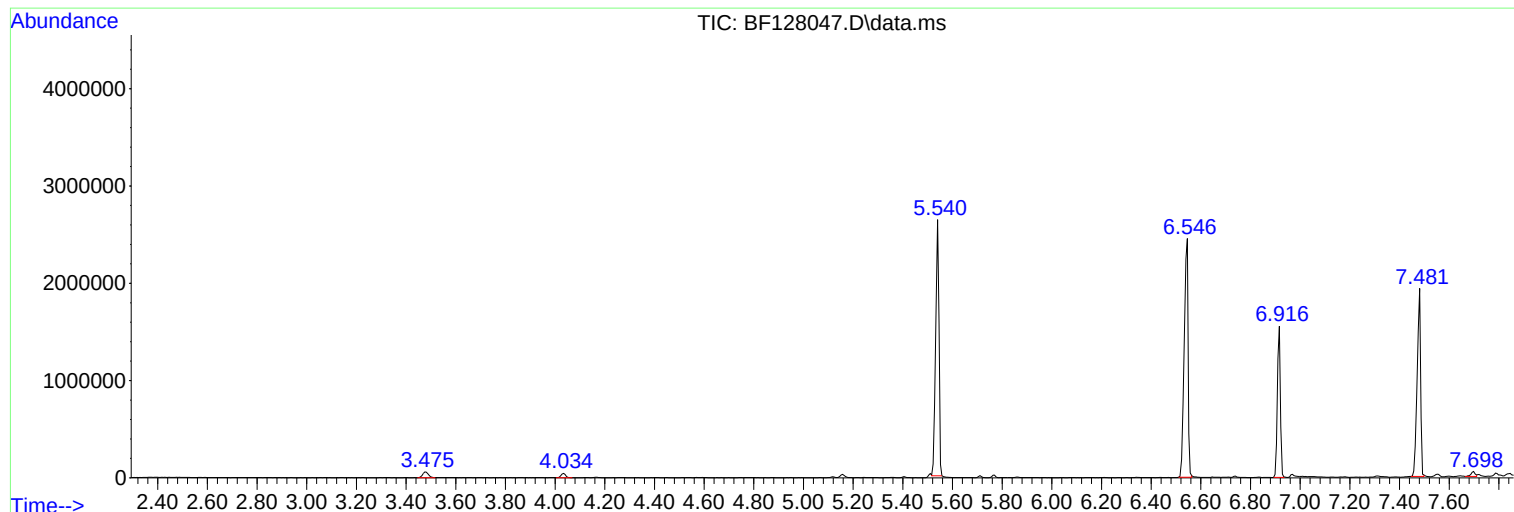
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Acq On : 03 May 2022 15:52
Operator : CG\JU
Sample : N2650-03
Misc :
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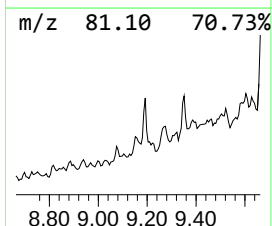
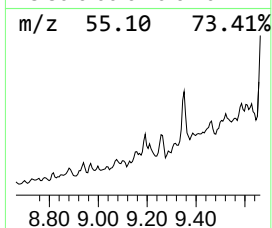
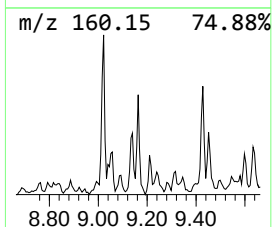
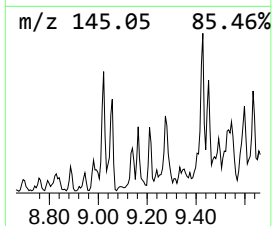
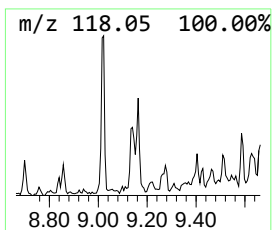
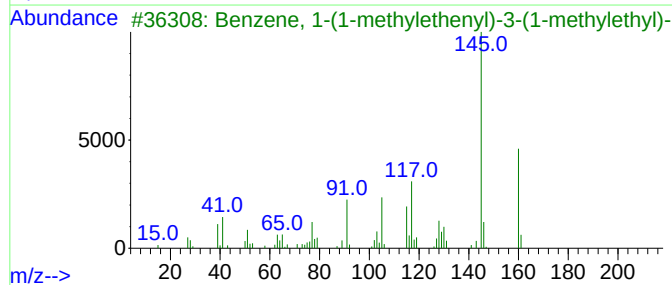
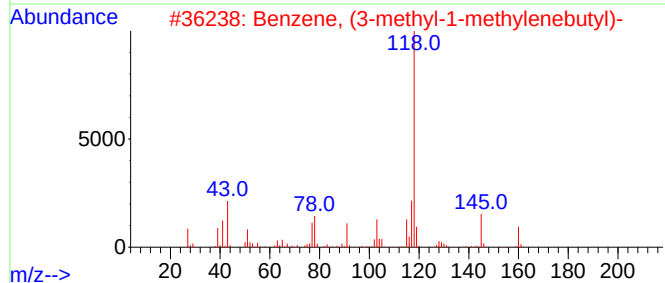
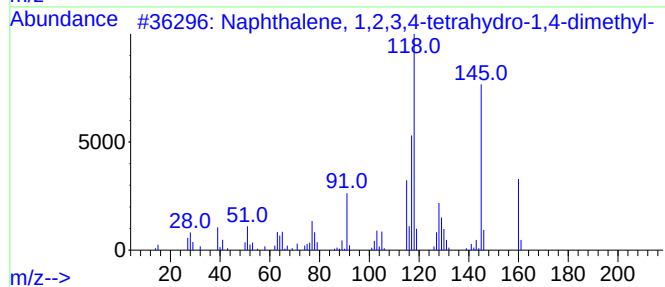
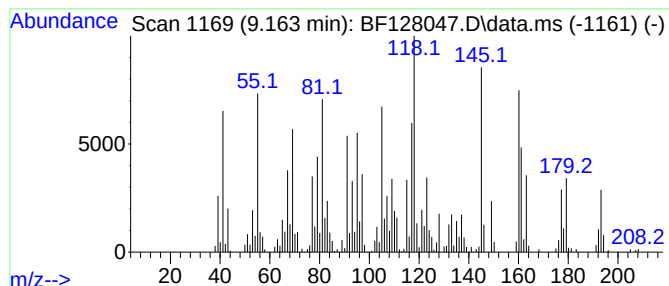
Instrument :
BNA_F
ClientSampleId :
ETGI-362

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
9.163	7.68 ng	597915	Acenaphthene-d10		9.969	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...		160	C12H16	004175-54-6	64
2	Benzene, (3-methyl-1-methylenebu...		160	C12H16	038212-14-5	53
3	Benzene, 1-(1-methylethenyl)-3-(...		160	C12H16	001129-29-9	46
4	Benzene, (1,3-dimethyl-2-butenyl)-		160	C12H16	050704-01-3	46
5	1H-Inden-1-one, 2,3-dihydro-3,3-...		160	C11H12O	026465-81-6	38



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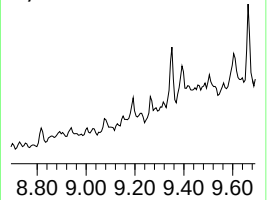
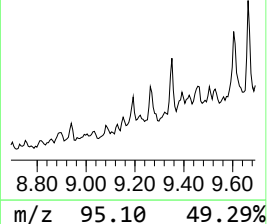
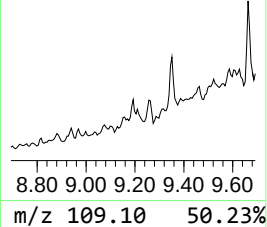
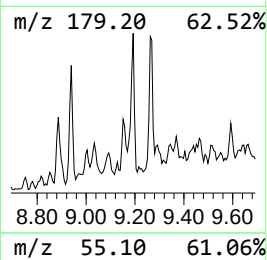
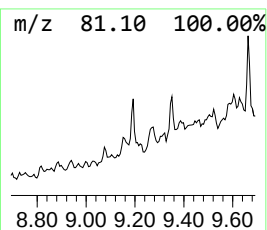
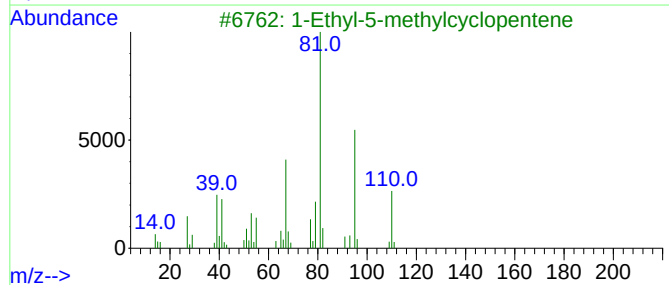
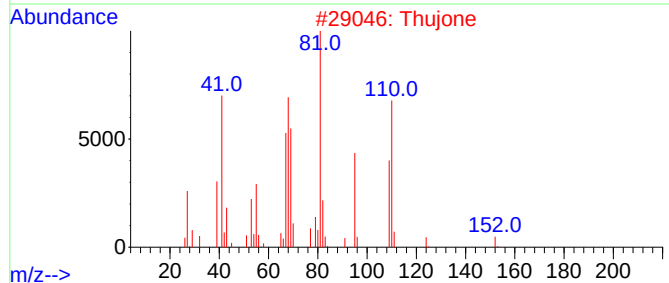
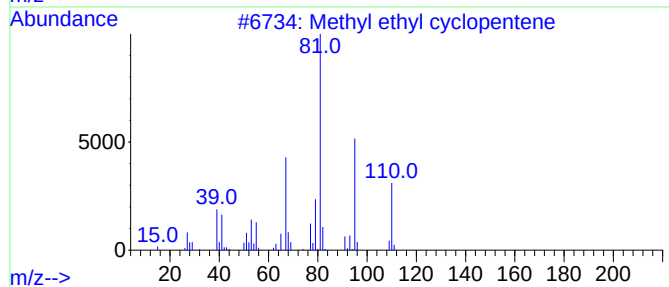
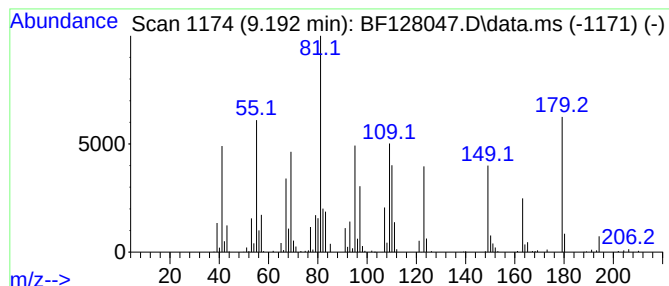
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown9.192 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.192	2.66 ng	207150	Acenaphthene-d10	9.969

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl ethyl cyclopentene	110	C8H14	019780-56-4	42
2			Thujone	152	C10H16O	000546-80-5	38
3			1-Ethyl-5-methylcyclopentene	110	C8H14	097797-57-4	27
4			2-Propenyl-3-vinyloxirane	110	C7H10O	070597-49-8	27
5			Cyclohexene, 1-ethyl-	110	C8H14	001453-24-3	25



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Misc :
ALS Vial : 9 Sample Multiplier: 1

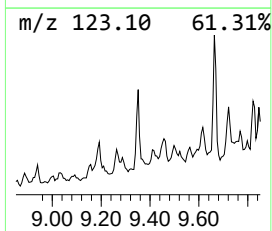
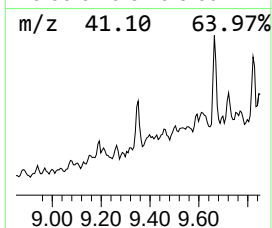
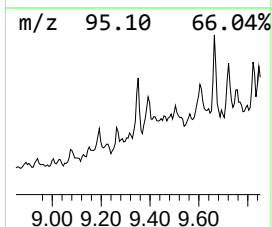
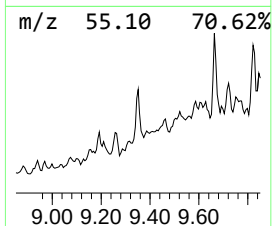
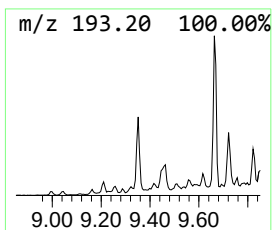
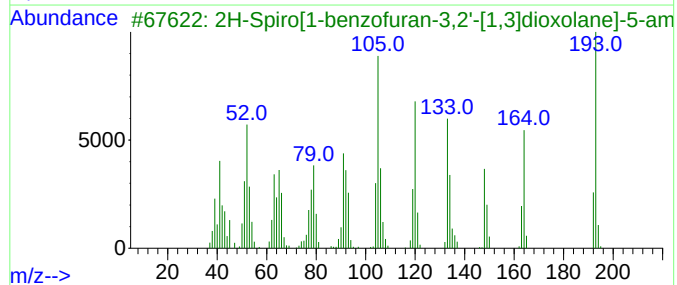
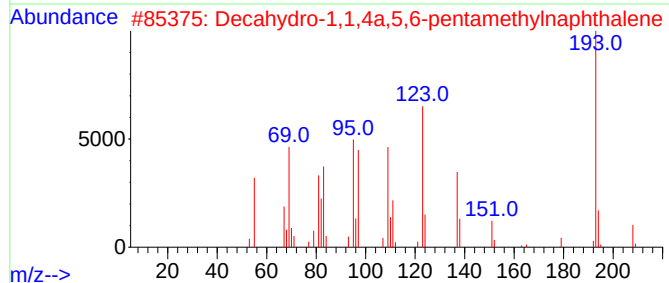
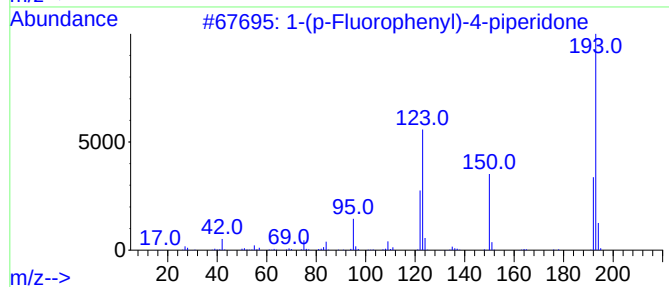
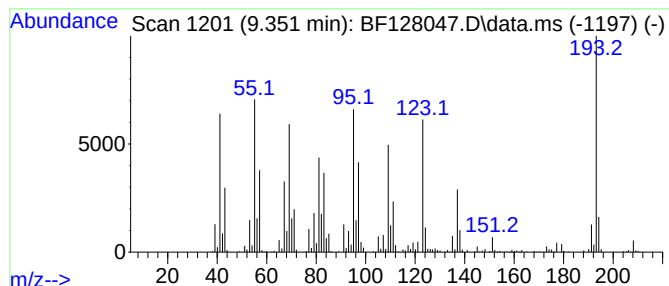
Instrument :
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 unknown9.351 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD			R.T.
9.351	10.98 ng	855165	Acenaphthene-d10			9.969
Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-(p-Fluorophenyl)-4-piperidone	193	C11H12FNO	1000238-56-7	43
2		Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	43
3		2H-Spiro[1-benzofuran-3,2'-[1,3]...	193	C10H11NO3	1000387-33-4	35
4		4-Amino-N-hydroxypyrazolo[3,2-c]...	193	C6H7N7O	1000443-50-2	27
5		9-Phenanthrenamine	193	C14H11N	000947-73-9	25



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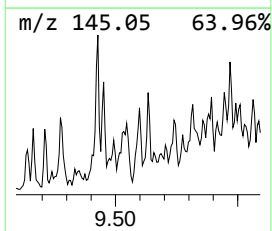
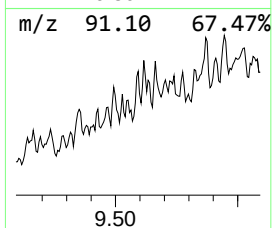
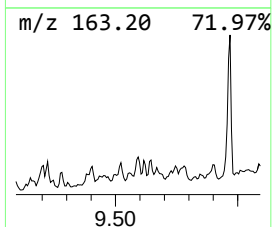
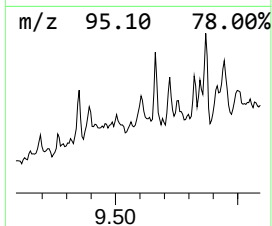
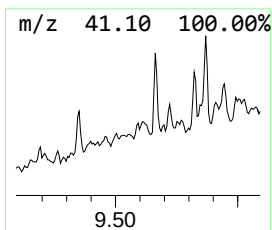
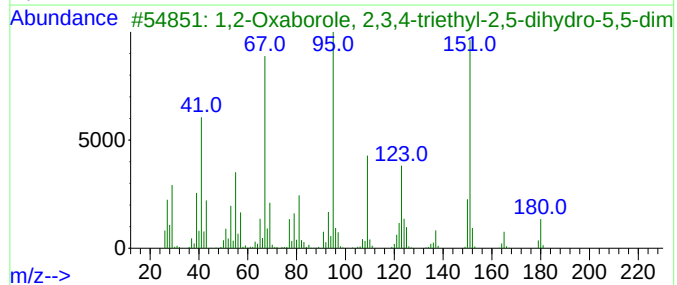
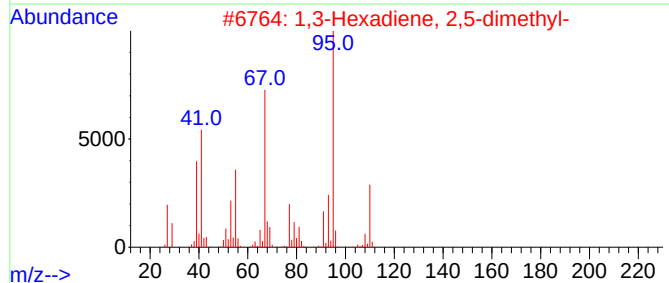
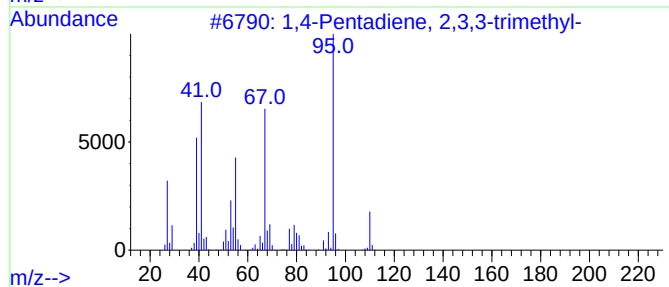
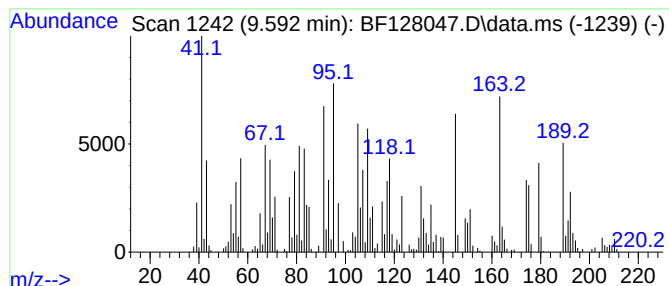
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 unknown9.592 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.592	3.68 ng	286168	Acenaphthene-d10	9.969

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1,4-Pentadiene, 2,3,3-trimethyl-	110	C8H14	000756-02-5	25
2		1,3-Hexadiene, 2,5-dimethyl-	110	C8H14	029253-64-3	11
3		1,2-Oxaborole, 2,3,4-triethyl-2,...	180	C11H21BO	061142-64-1	11
4		1,4-Dichloro-2-nitrosobenzene	175	C6H3Cl2NO	067083-41-4	11
5		Cyclopentane, 1,2-dimethyl-3-met...	110	C8H14	1000150-99-3	11



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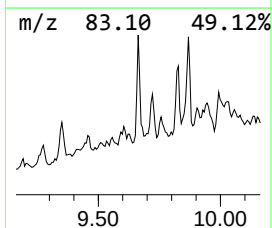
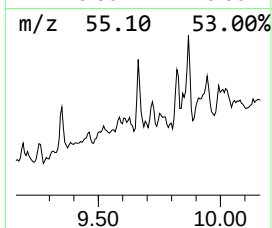
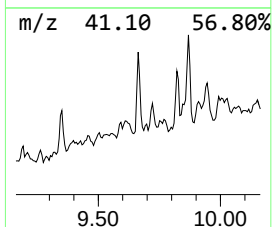
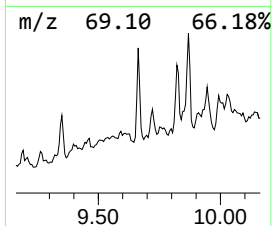
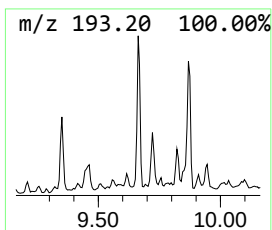
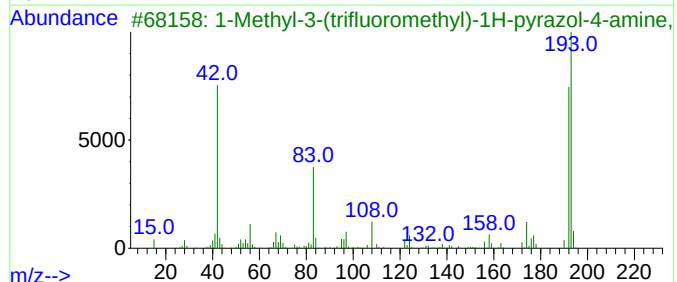
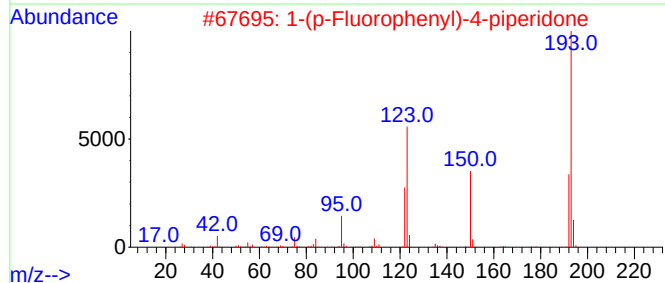
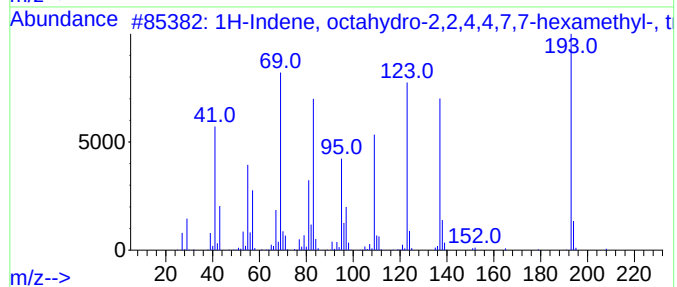
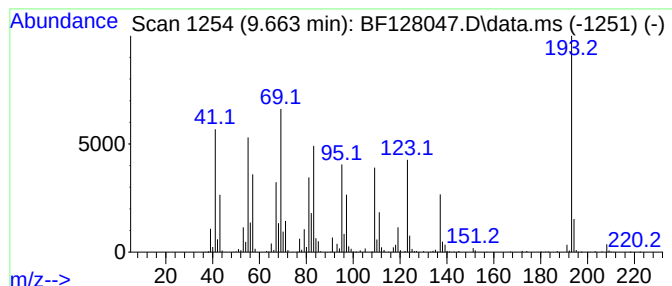
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 unknown9.663 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.663	15.10 ng	1175990	Acenaphthene-d10	9.969

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, octahydro-2,2,4,7,7...	208	C15H28	054832-83-6	40		
2	1-(p-Fluorophenyl)-4-piperidone	193	C11H12FNO	1000238-56-7	35		
3	1-Methyl-3-(trifluoromethyl)-1H...	193	C7H10F3N3	1000511-73-1	35		
4	9-Phenanthrenamine	193	C14H11N	000947-73-9	22		
5	2H-Spiro[1-benzofuran-3,2'-[1,3]...	193	C10H11NO3	1000387-33-4	18		



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050322\
Data File : BF128047.D
Acq On : 03 May 2022 15:52
Operator : CG\JU
Sample : N2650-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ETGI-362

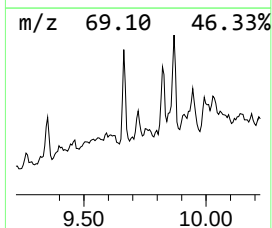
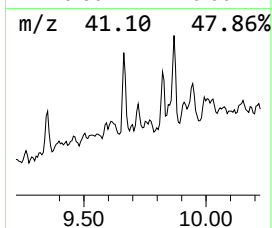
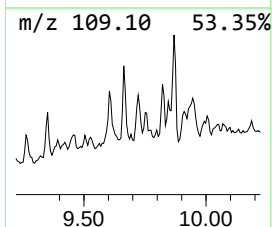
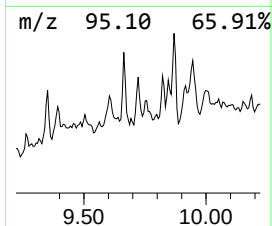
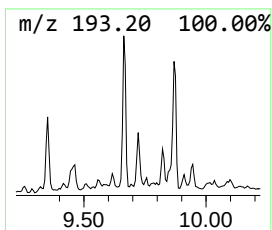
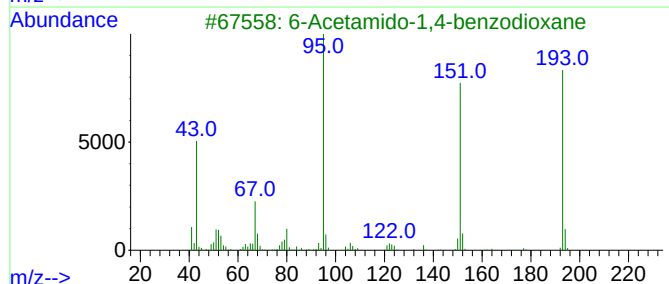
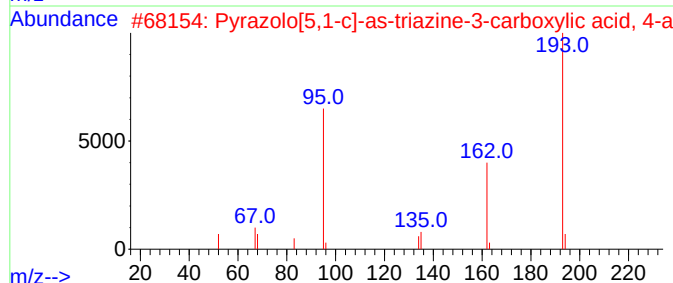
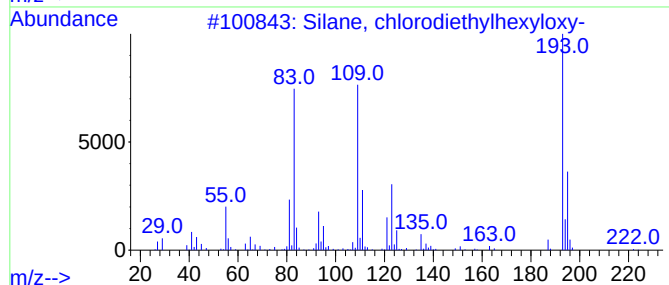
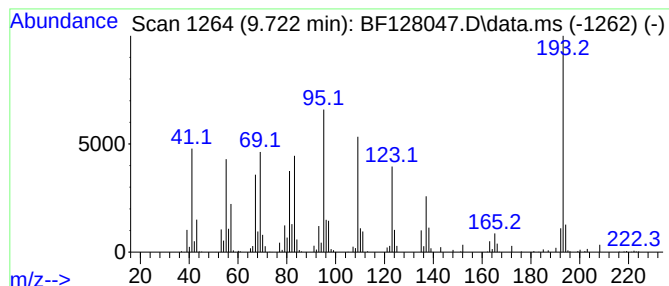
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Silane, chlorodiethylhexyloxy- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.722	5.26 ng	409356	Acenaphthene-d10	9.969

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Silane, chlorodiethylhexyloxy-	222	C10H23ClOSi	1000363-74-4	52
2			Pyrazolo[5,1-c]-as-triazine-3-ca...	193	C6H7N7O	016111-78-7	35
3			6-Acetamido-1,4-benzodioxane	193	C10H11NO3	063546-19-0	35
4			Cyclopentene, 1-isopropyl-2,3-di...	138	C10H18	007712-73-4	35
5			3-Phenylindole	193	C14H11N	001504-16-1	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050322\
Data File : BF128047.D
Acq On : 03 May 2022 15:52
Operator : CG\JU
Sample : N2650-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ETGI-362

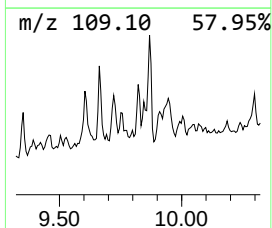
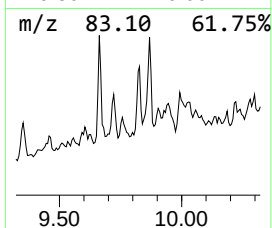
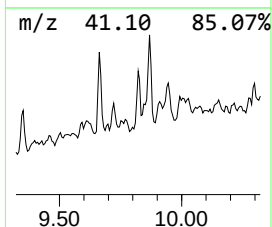
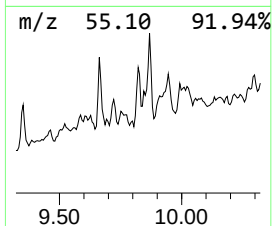
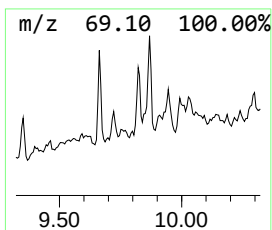
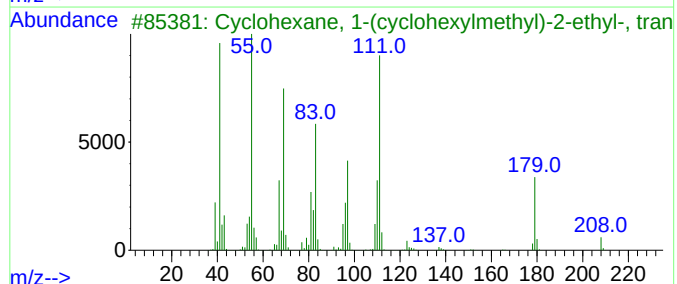
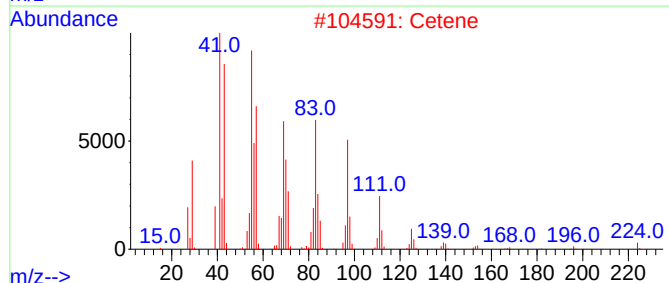
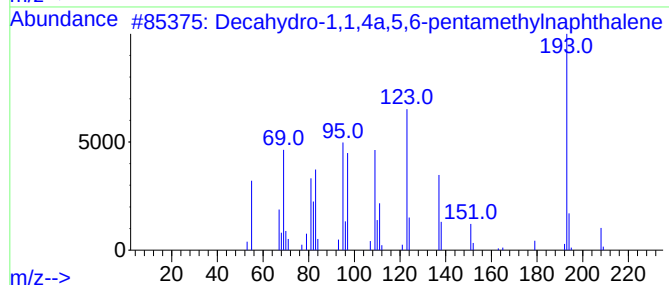
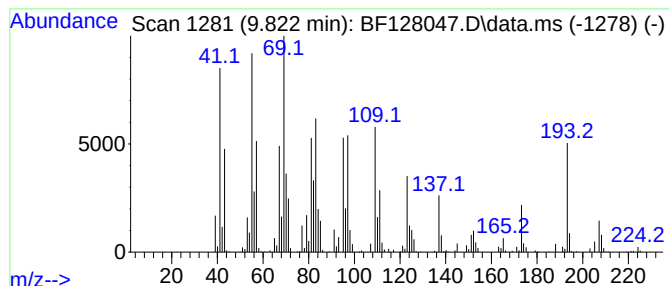
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Decahydro-1,1,4a,5,6-pentam... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.822	11.21 ng	873012	Acenaphthene-d10	9.969

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	70
2			Cetene	224	C16H32	000629-73-2	40
3			Cyclohexane, 1-(cyclohexylmethyl)...	208	C15H28	054934-92-8	38
4			Silane, chlorodiethyl(2-methylpe...	222	C10H23ClOSi	1000363-79-1	38
5			Cyclopropane carboxamide, 2-cycl...	207	C13H21NO	331416-19-4	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050322\
Data File : BF128047.D
Acq On : 03 May 2022 15:52
Operator : CG\JU
Sample : N2650-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

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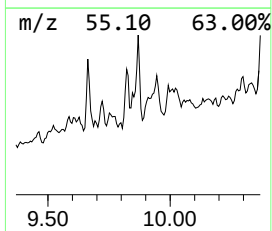
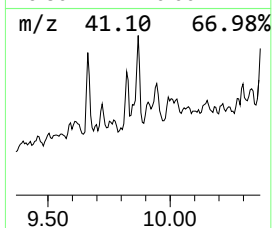
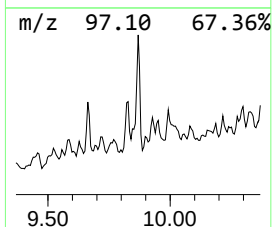
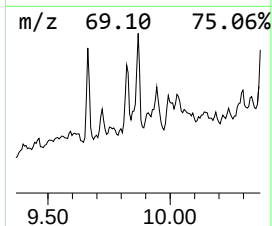
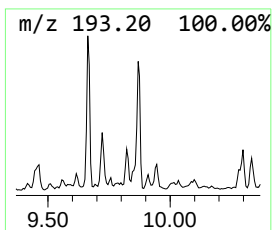
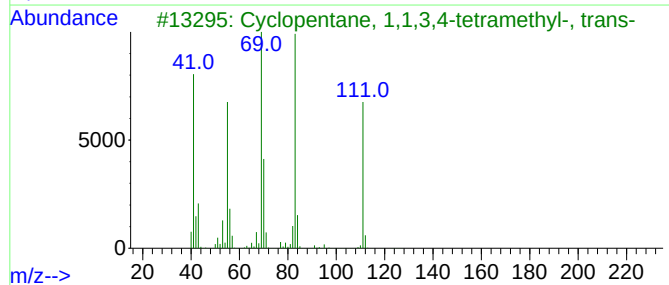
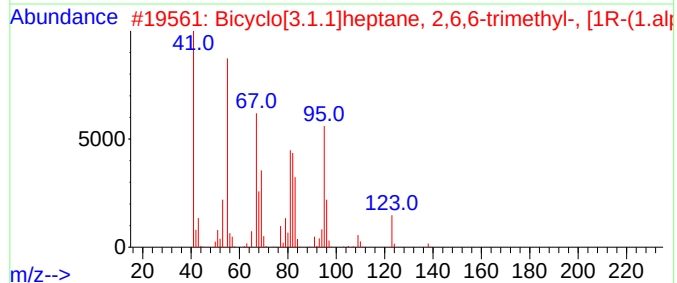
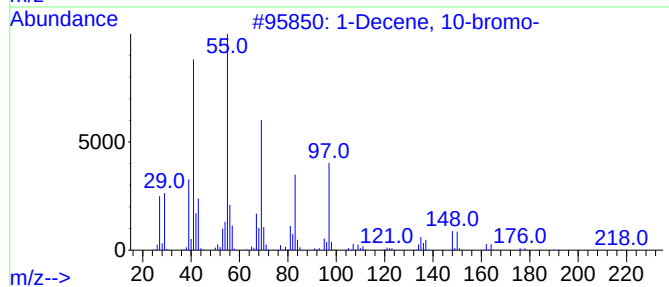
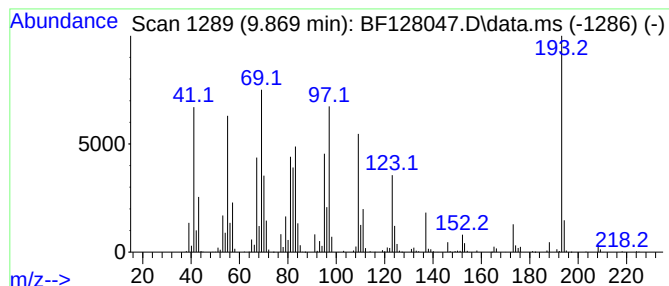
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown9.869 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.
9.869	17.62 ng	1371960	Acenaphthene-d10		9.969

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Decene, 10-bromo-	218	C10H19Br	062871-09-4	27
2		Bicyclo[3.1.1]heptane, 2,6,6-tri...	138	C10H18	004795-86-2	25
3		Cyclopentane, 1,1,3,4-tetramethy...	126	C9H18	020309-77-7	22
4		Oxiranecarboxaldehyde, 3-methyl-...	168	C10H16O2	016996-12-6	22
5		1,6-Heptadiene, 2,5,5-trimethyl-	138	C10H18	062238-28-2	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050322\
Data File : BF128047.D
Acq On : 03 May 2022 15:52
Operator : CG\JU
Sample : N2650-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

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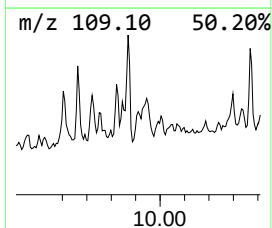
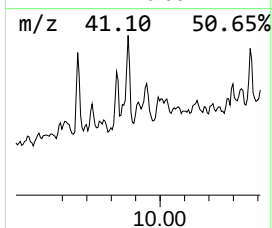
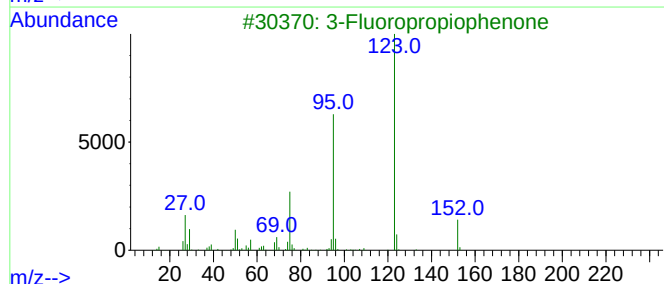
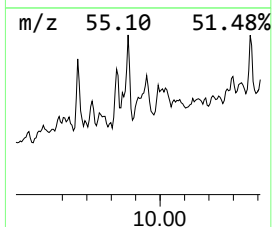
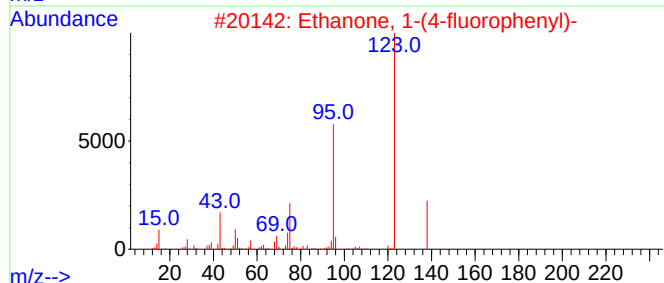
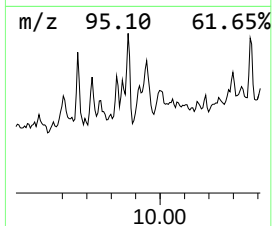
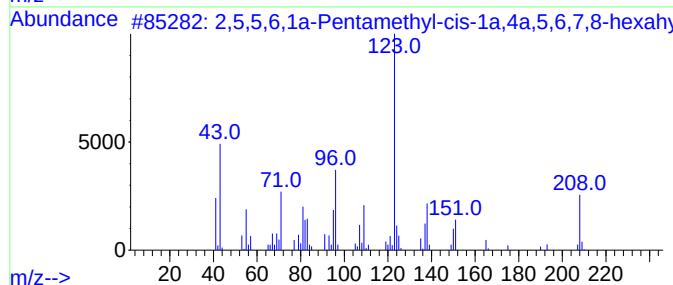
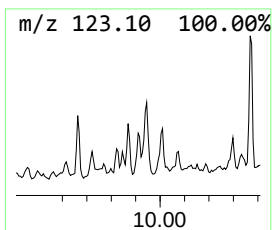
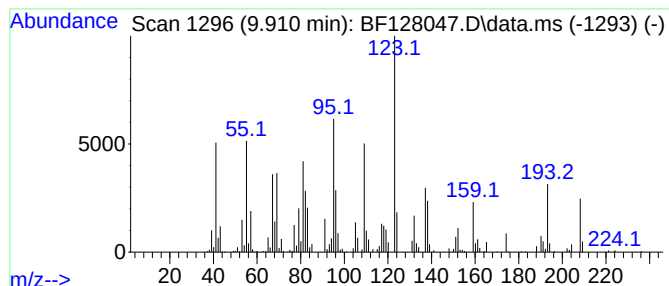
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown9.910 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.910	6.48 ng	504463	Acenaphthene-d10	9.969

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,5,5,6,1a-Pentamethyl-cis-1a,4a...	208	C14H24O	1010215-77-9	49		
2	Ethanone, 1-(4-fluorophenyl)-	138	C8H7FO	000403-42-9	43		
3	3-Fluoropropiophenone	152	C9H9FO	000455-67-4	27		
4	1-(2,2,6-Trimethylcyclohexyl)hex...	268	C17H32O2	1000498-96-1	27		
5	o-Isobutyranisidide	193	C11H15NO2	071182-38-2	25		



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050322\
Data File : BF128047.D
Acq On : 03 May 2022 15:52
Operator : CG\JU
Sample : N2650-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

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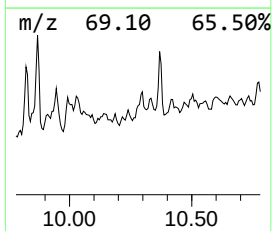
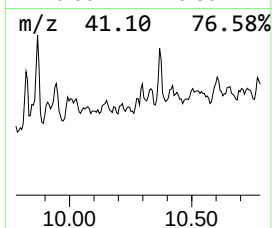
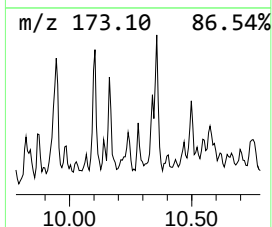
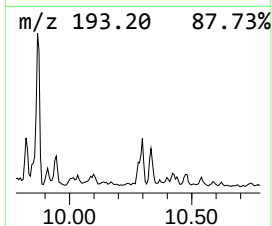
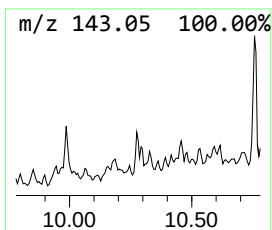
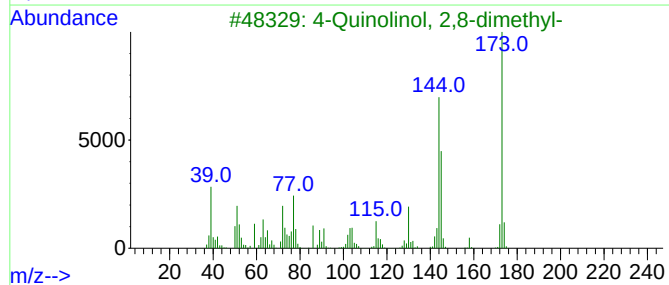
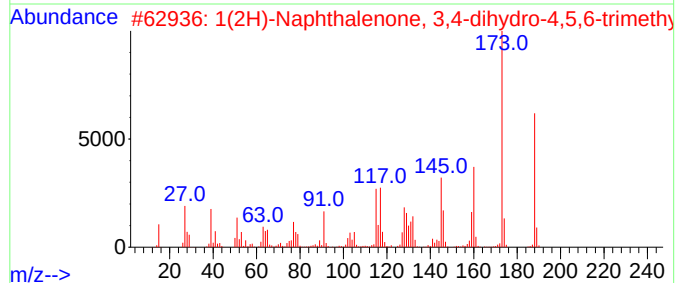
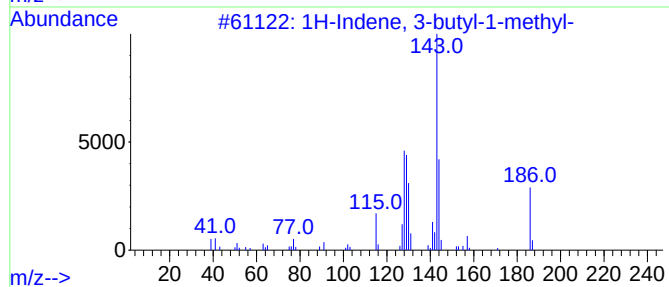
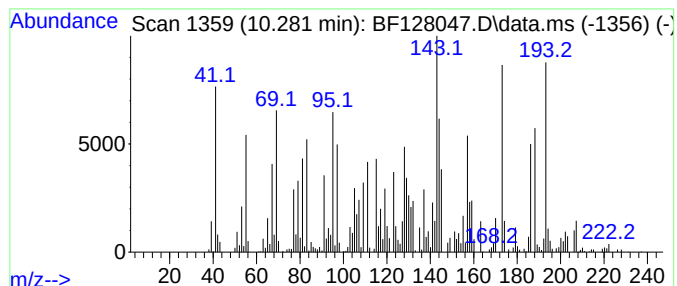
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown10.281 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.281	5.21 ng	405493	Acenaphthene-d10	9.969

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 3-butyl-1-methyl-	186	C14H18	111400-84-1	25
2			1(2H)-Naphthalenone, 3,4-dihydro...	188	C13H16O	030316-31-5	25
3			4-Quinolinol, 2,8-dimethyl-	173	C11H11NO	1010351-37-0	20
4			1,1,4,5,6-Pentamethyl-2,3-dihydr...	188	C14H20	016204-67-4	18
5			(1-Methylenepent-2-enyl)benzene	158	C12H14	084313-24-6	15



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050322\
Data File : BF128047.D
Acq On : 03 May 2022 15:52
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Sample : N2650-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

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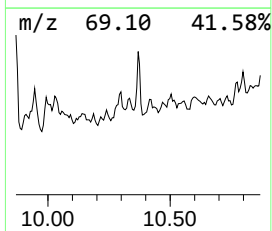
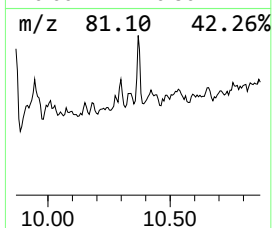
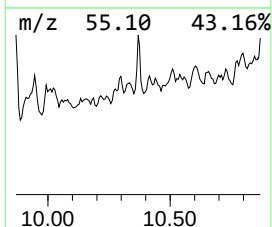
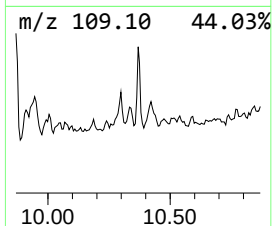
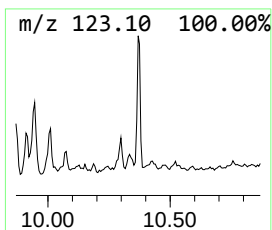
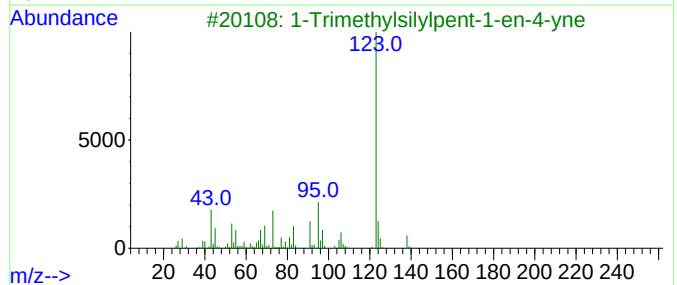
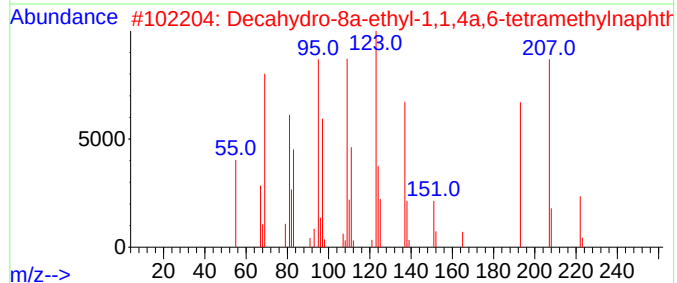
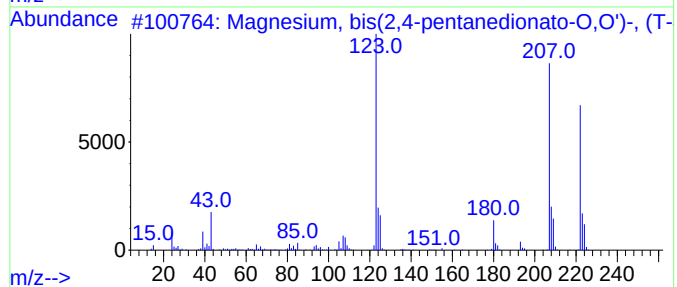
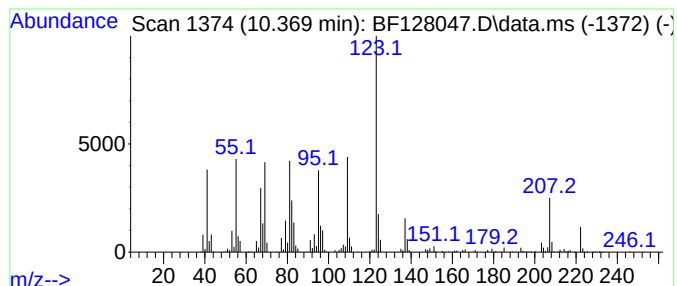
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown10.369 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.369	9.07 ng	705973	Acenaphthene-d10	9.969

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Magnesium, bis(2,4-pentanedionat...	222	C10H14MgO4	014024-56-7	47
2			Decahydro-8a-ethyl-1,1,4a,6-tetr...	222	C16H30	1000100-23-6	46
3			1-Trimethylsilylpent-1-en-4-yne	138	C8H14Si	079516-25-9	43
4			1,3-Benzenediamine, 4-methoxy-	138	C7H10N2O	000615-05-4	43
5			Benzamide, 4-fluoro-N-(2,6-dimet...	243	C15H14FNO	117805-18-2	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050322\
Data File : BF128047.D
Acq On : 03 May 2022 15:52
Operator : CG\JU
Sample : N2650-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ETGI-362

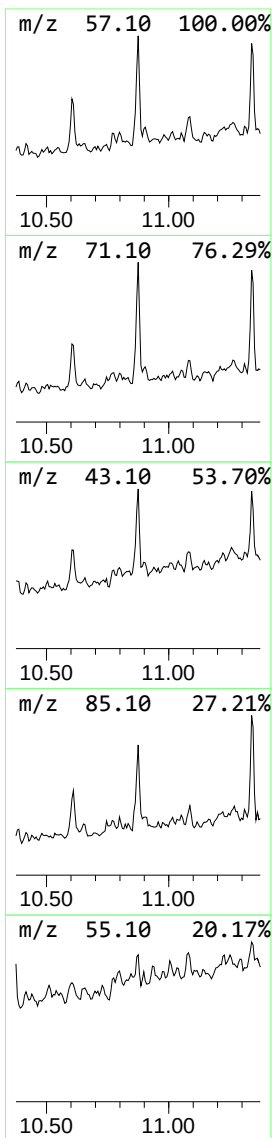
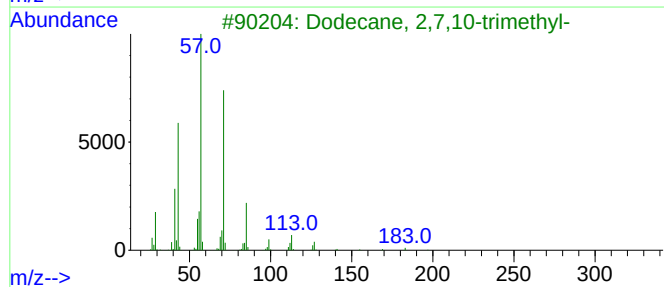
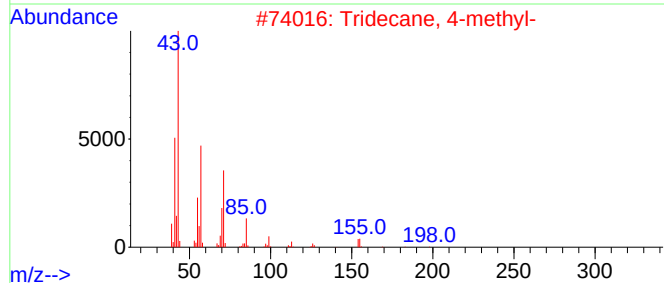
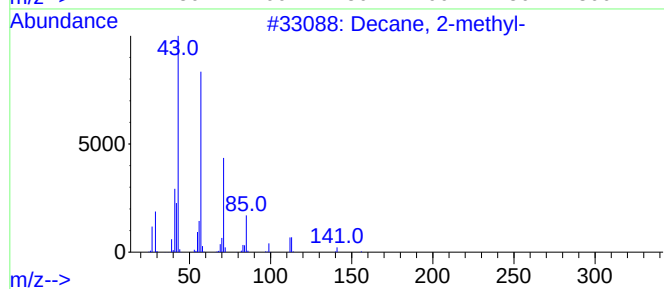
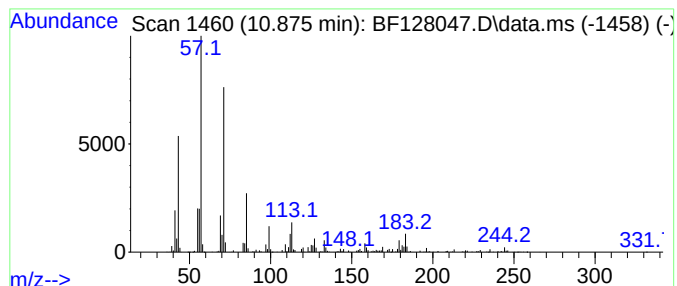
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Decane, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.875	7.15 ng	538127	Phenanthrene-d10	11.463

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decane, 2-methyl-	156	C11H24	006975-98-0	87
2		Tridecane, 4-methyl-	198	C14H30	026730-12-1	86
3		Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	80
4		Tridecane, 5-propyl-	226	C16H34	055045-11-9	80
5		Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050322\
Data File : BF128047.D
Acq On : 03 May 2022 15:52
Operator : CG\JU
Sample : N2650-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

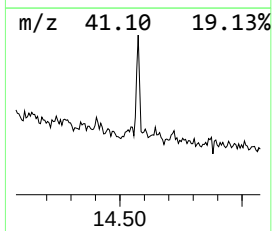
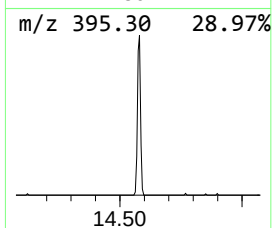
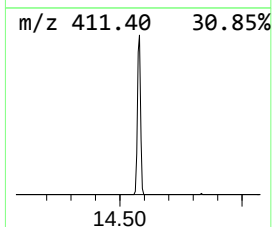
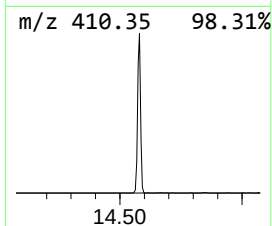
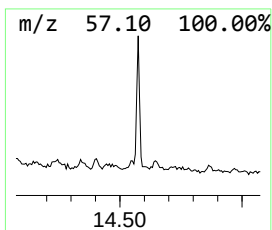
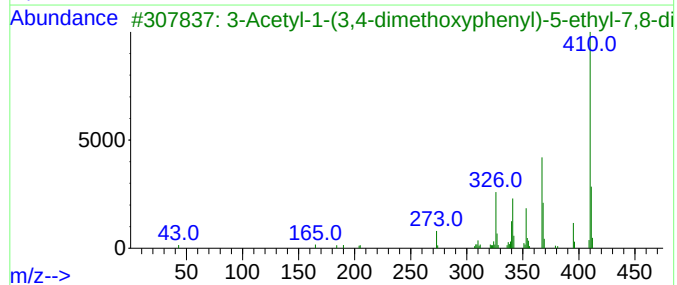
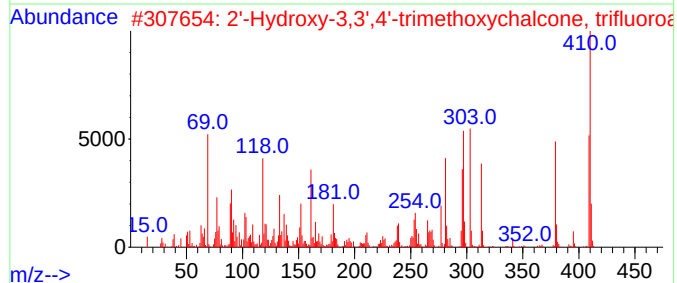
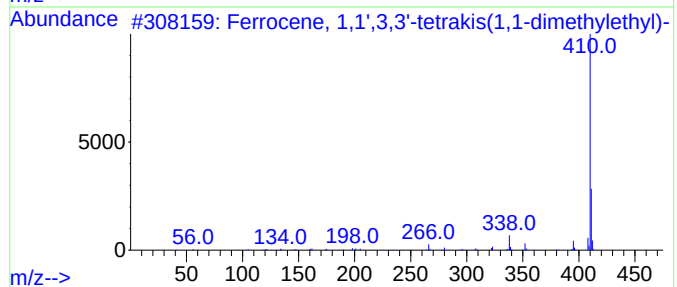
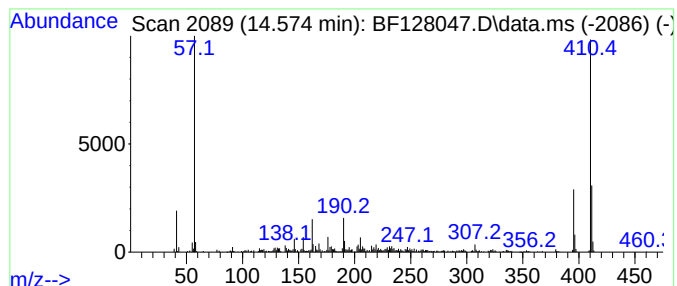
Instrument :
BNA_F
ClientSampleId :
ETGI-362

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Ferrocene, 1,1',3,3'-tetrak... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD			R.T.
14.574	6.04 ng	305903	Chrysene-d12			14.104
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Ferrocene, 1,1',3,3'-tetrakis(1,...		410	C26H42Fe	001317-17-5	60
2	2'-Hydroxy-3,3',4'-trimethoxycha...		410	C20H17F3O6	1010453-84-8	46
3	3-Acetyl-1-(3,4-dimethoxyphenyl)...		410	C23H26N2O5	090140-65-1	40
4	6,9-Methanobenzo[h]pentaphene-5,...		410	C27H22O4	057427-50-6	38
5	9-Chloro-4,5-dihydro-4-[3,4,5-tr...		410	C22H19ClN2O4	1000212-48-1	33



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050322\
Data File : BF128047.D
Acq On : 03 May 2022 15:52
Operator : CG\JU
Sample : N2650-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ETGI-362

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042222.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
Naphthalene, 1,...	9.163	7.7	ng	597915	3	9.969	1557170	20.0
unknown9.192	9.192	2.7	ng	207150	3	9.969	1557170	20.0
unknown9.351	9.351	11.0	ng	855165	3	9.969	1557170	20.0
unknown9.592	9.592	3.7	ng	286168	3	9.969	1557170	20.0
unknown9.663	9.663	15.1	ng	1175990	3	9.969	1557170	20.0
Silane, chlorod...	9.722	5.3	ng	409356	3	9.969	1557170	20.0
Decahydro-1,1,4...	9.822	11.2	ng	873012	3	9.969	1557170	20.0
unknown9.869	9.869	17.6	ng	1371960	3	9.969	1557170	20.0
unknown9.910	9.910	6.5	ng	504463	3	9.969	1557170	20.0
unknown10.281	10.281	5.2	ng	405493	3	9.969	1557170	20.0
unknown10.369	10.369	9.1	ng	705973	3	9.969	1557170	20.0
Decane, 2-methyl-	10.875	7.2	ng	538127	4	11.463	1506000	20.0
Ferrocene, 1,1'...	14.574	6.0	ng	305903	5	14.104	1012970	20.0