

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112025\
 Data File : BF144299.D
 Acq On : 20 Nov 2025 17:27
 Operator : RC/JU
 Sample : Q3681-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ETGI-354

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-BF110525.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF144299.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.081	483	491	501	rBV	205555	303840	26.37%	2.926%
2	5.487	555	560	582	rBV	735280	1049687	91.11%	10.107%
3	6.498	726	732	753	rBV	739426	1063264	92.29%	10.238%
4	6.863	789	794	799	rBV	197182	238123	20.67%	2.293%
5	7.422	878	889	899	rBV	449240	610280	52.97%	5.876%
6	7.881	961	967	974	rBV4	7009	11533	1.00%	0.111%
7	8.145	998	1012	1018	rBV	224900	311557	27.04%	3.000%
8	8.363	1040	1049	1054	rBV9	6965	15284	1.33%	0.147%
9	8.428	1054	1060	1064	rVB7	7059	13293	1.15%	0.128%
10	8.563	1080	1083	1088	rVV	9904	12978	1.13%	0.125%
11	8.857	1130	1133	1137	rVB	11232	13393	1.16%	0.129%
12	9.157	1179	1184	1189	rBV2	28277	43973	3.82%	0.423%
13	9.222	1189	1195	1200	rBV	771339	980507	85.10%	9.441%
14	9.416	1225	1228	1235	rVB	16168	27795	2.41%	0.268%
15	9.486	1235	1240	1245	rBV	44094	67286	5.84%	0.648%
16	9.557	1248	1252	1259	rVV	59061	97287	8.44%	0.937%
17	9.616	1259	1262	1268	rVV2	29265	43420	3.77%	0.418%
18	9.681	1268	1273	1282	rVB	25299	49180	4.27%	0.474%
19	9.763	1283	1287	1291	rBV	33106	40874	3.55%	0.394%
20	9.904	1302	1311	1315	rBV	228354	319272	27.71%	3.074%
21	9.933	1315	1316	1320	rVV2	27751	27936	2.42%	0.269%
22	10.004	1324	1328	1333	rVV2	23900	34617	3.00%	0.333%
23	10.051	1333	1336	1339	rVV3	8239	11732	1.02%	0.113%
24	10.104	1340	1345	1349	rVV2	40895	60749	5.27%	0.585%
25	10.145	1349	1352	1361	rVB2	31579	50481	4.38%	0.486%
26	10.222	1361	1365	1367	rBV	23073	33318	2.89%	0.321%
27	10.245	1367	1369	1375	rVB2	15629	19696	1.71%	0.190%
28	10.310	1375	1380	1386	rBV3	18447	42640	3.70%	0.411%
29	10.451	1399	1404	1410	rBV4	22544	44188	3.84%	0.425%
30	10.545	1417	1420	1426	rVB2	25608	39961	3.47%	0.385%
31	10.616	1427	1432	1437	rVB	105461	135770	11.78%	1.307%
32	10.692	1438	1445	1457	rBV	487854	652005	56.59%	6.278%
33	10.816	1461	1466	1473	rVB	47889	76191	6.61%	0.734%
34	10.928	1481	1485	1489	rVB	16009	18617	1.62%	0.179%
35	10.998	1492	1497	1499	rVV4	11415	19459	1.69%	0.187%
36	11.028	1499	1502	1514	rVB4	21409	51506	4.47%	0.496%

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Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

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37	11.145	1514	1522	1527	rBV6	12517	36492	3.17%	0.351%
38	11.245	1535	1539	1541	rBV4	8795	13957	1.21%	0.134%
39	11.280	1542	1545	1551	rVB2	23351	38811	3.37%	0.374%
40	11.392	1559	1564	1574	rBV2	253674	464495	40.32%	4.472%
41	11.469	1574	1577	1580	rVV	17243	20706	1.80%	0.199%
42	11.504	1580	1583	1588	rVB4	8227	12786	1.11%	0.123%
43	11.627	1599	1604	1610	rBV3	17371	32635	2.83%	0.314%
44	11.916	1646	1653	1659	rBV2	64319	150613	13.07%	1.450%
45	12.010	1665	1669	1676	rVB	33993	64388	5.59%	0.620%
46	12.445	1741	1743	1746	rBV	13825	15502	1.35%	0.149%
47	12.610	1766	1771	1776	rBV	224577	295994	25.69%	2.850%
48	12.839	1804	1810	1816	rBV	190620	279025	24.22%	2.687%
49	12.980	1829	1834	1840	rVB	903608	1152151	100.00%	11.094%
50	13.233	1875	1877	1880	rBV	19687	21875	1.90%	0.211%
51	13.880	1984	1987	1993	rVB3	67768	107984	9.37%	1.040%
52	14.039	2009	2014	2018	rBV2	319845	543497	47.17%	5.233%
53	15.092	2189	2193	2200	rBV	85423	168719	14.64%	1.625%
54	15.521	2262	2266	2279	rVB	173728	334432	29.03%	3.220%

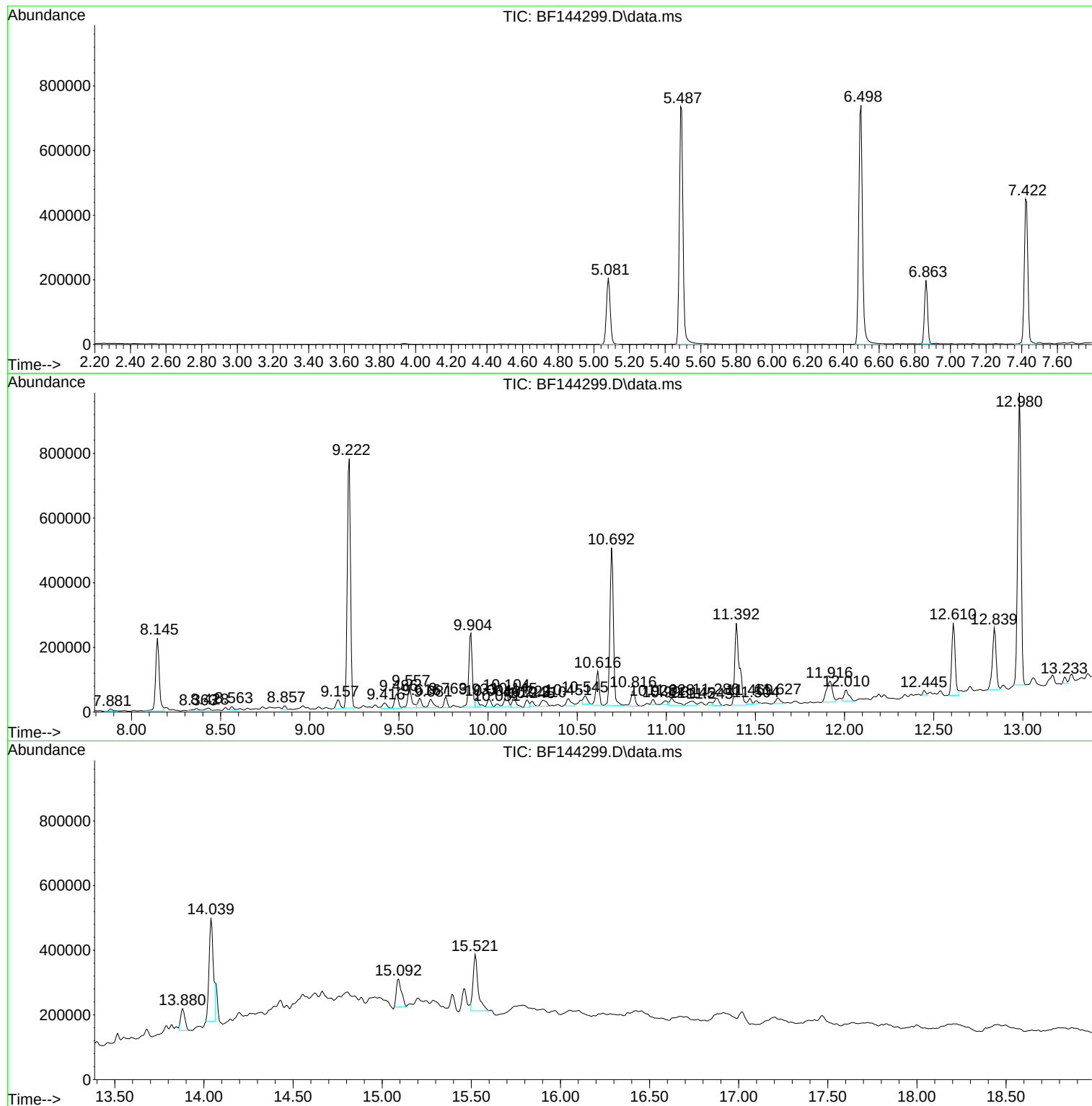
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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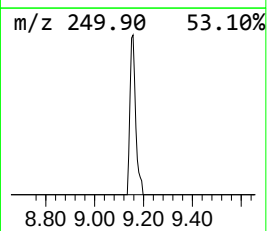
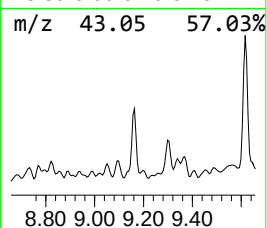
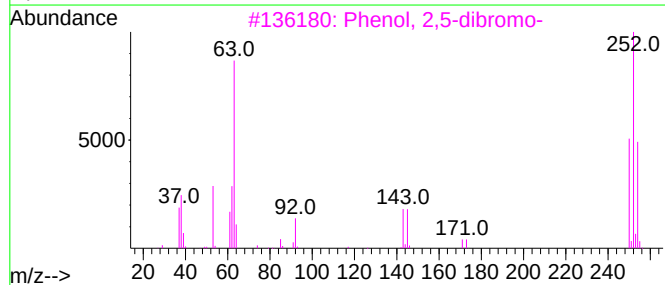
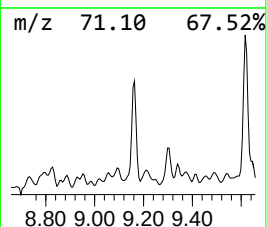
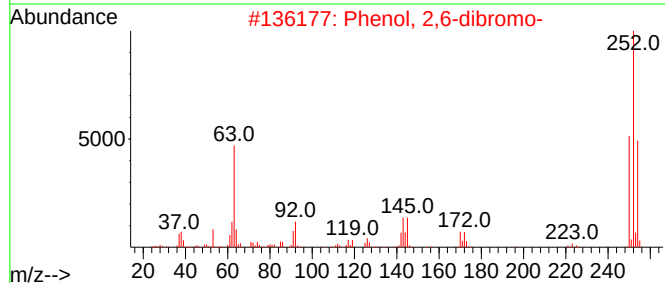
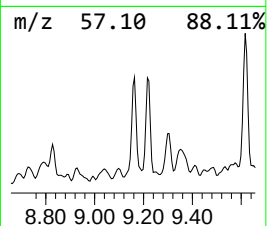
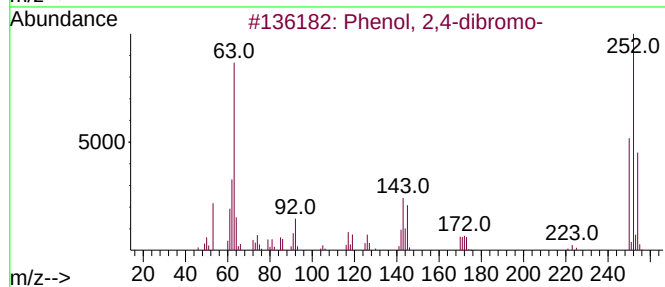
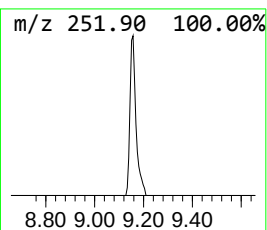
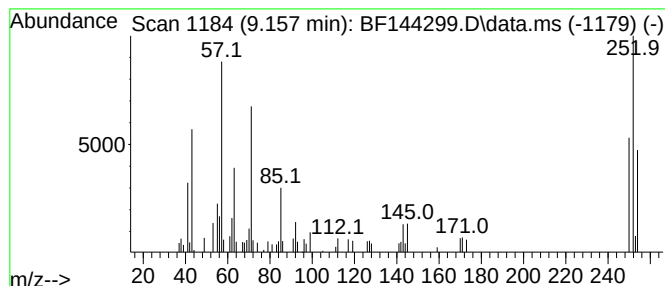
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Phenol, 2,4-dibromo- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.157	2.75 ng	43973	Acenaphthene-d10	9.904

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,4-dibromo-	250	C6H4Br2O	000615-58-7	96
2			Phenol, 2,6-dibromo-	250	C6H4Br2O	000608-33-3	96
3			Phenol, 2,5-dibromo-	250	C6H4Br2O	028165-52-8	95
4			3,4-Dibromophenol	250	C6H4Br2O	1010487-11-5	91
5			Phenol, 2,3-dibromo-	250	C6H4Br2O	057383-80-9	87



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Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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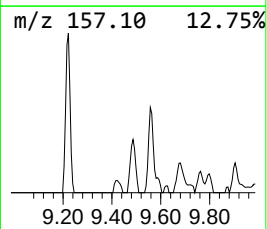
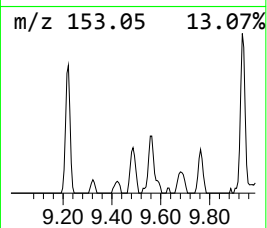
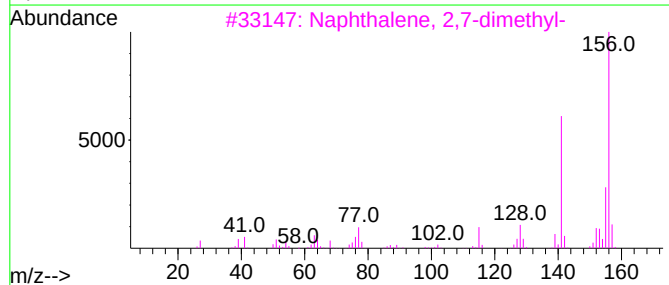
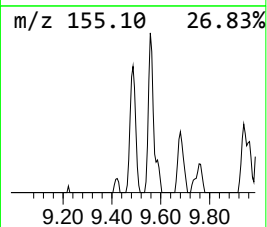
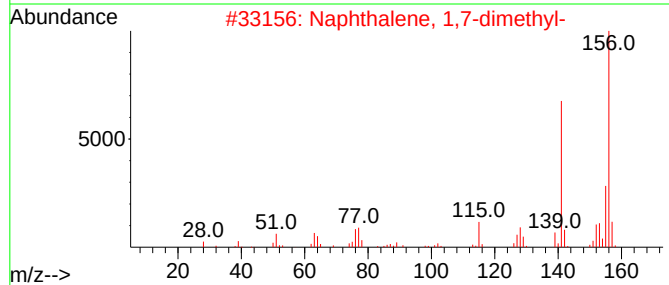
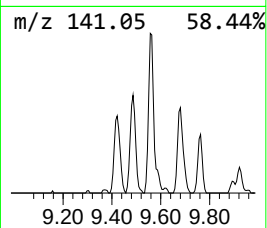
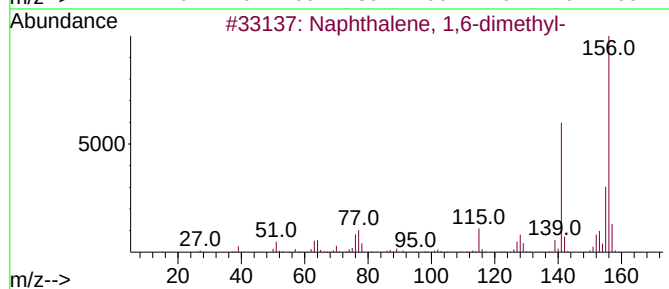
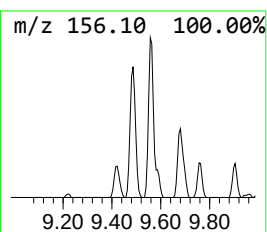
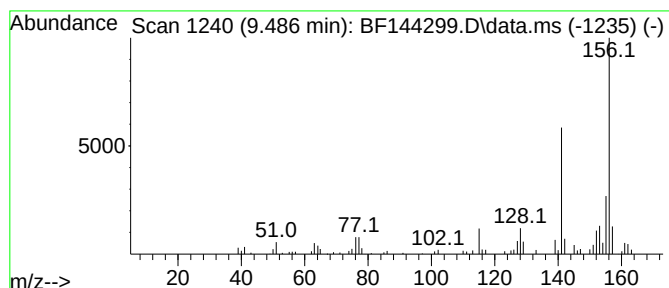
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Naphthalene, 1,6-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.486	4.21 ng	67286	Acenaphthene-d10	9.904

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
2			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
3			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
4			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96
5			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	96



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Misc :
ALS Vial : 11 Sample Multiplier: 1

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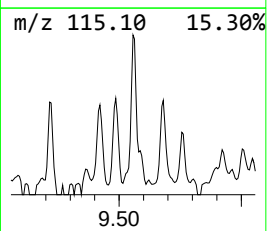
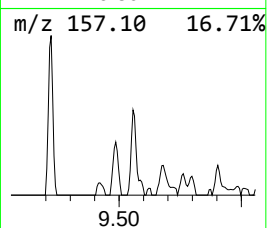
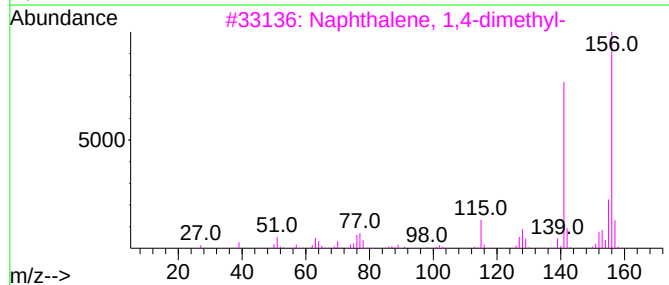
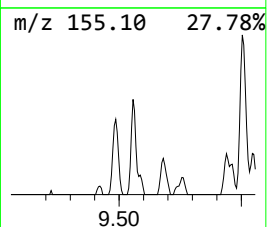
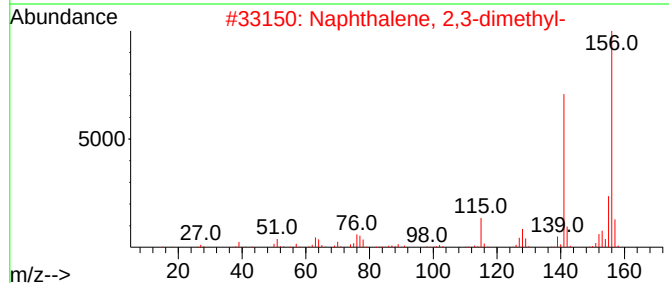
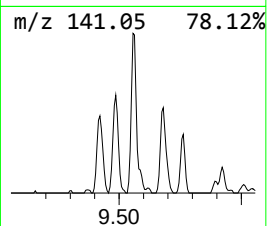
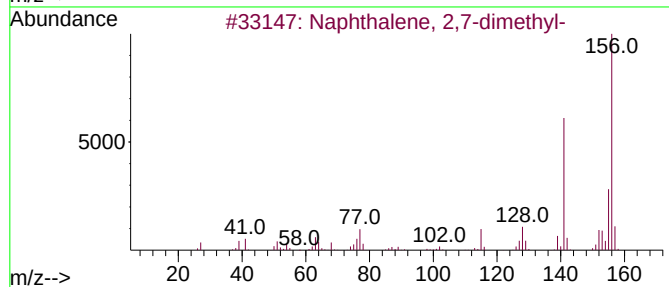
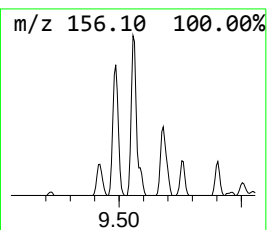
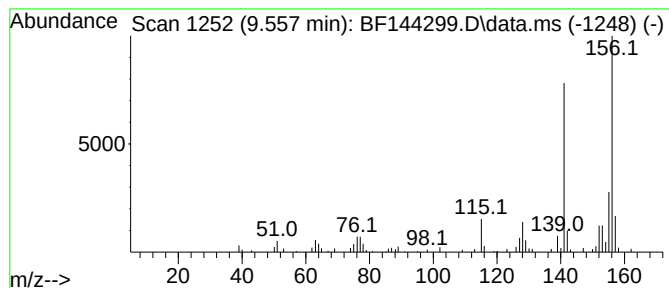
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Naphthalene, 2,7-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.557	6.09 ng	97287	Acenaphthene-d10	9.904

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



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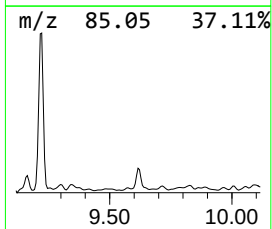
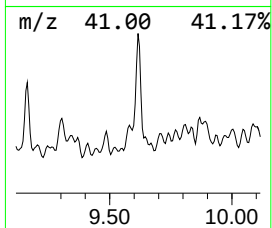
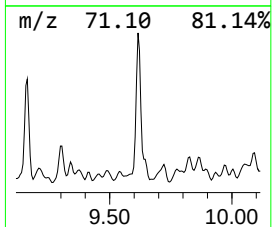
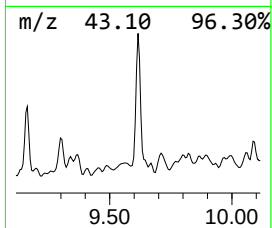
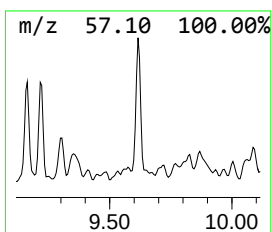
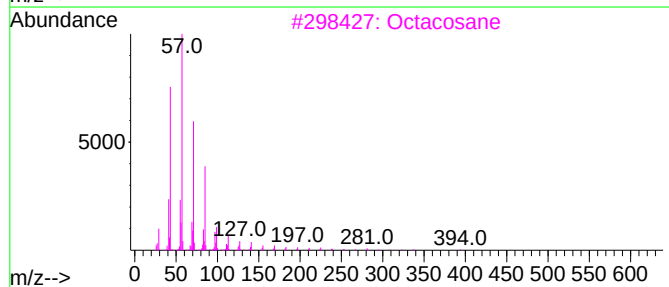
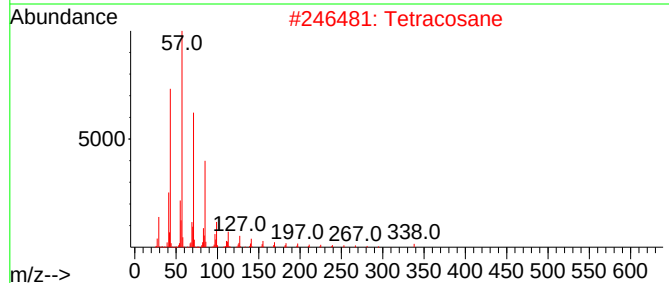
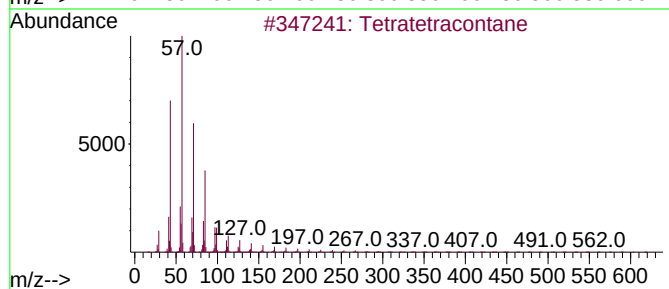
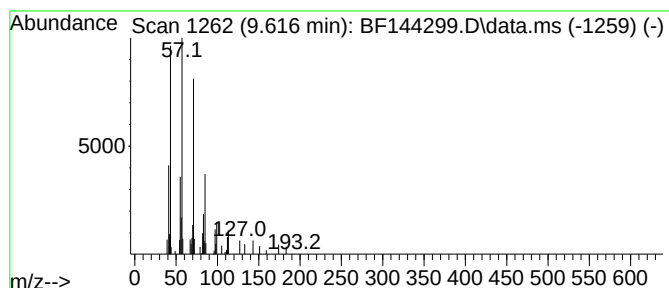
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Tetratetracontane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.616	2.72 ng	43420	Acenaphthene-d10	9.904

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetratetracontane	619	C44H90	007098-22-8	86
2			Tetracosane	338	C24H50	000646-31-1	86
3			Octacosane	394	C28H58	000630-02-4	86
4			Hexatriacontane	507	C36H74	000630-06-8	80
5			2-Bromo dodecane	248	C12H25Br	013187-99-0	80



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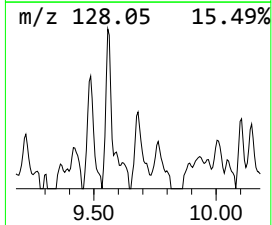
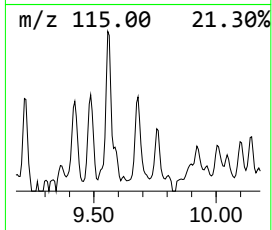
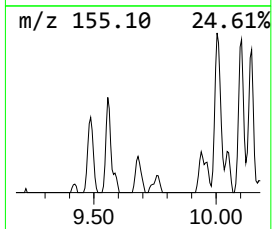
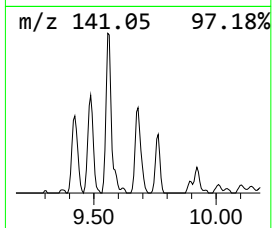
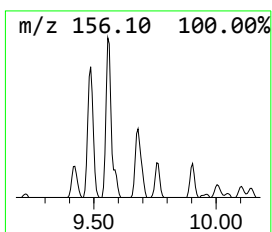
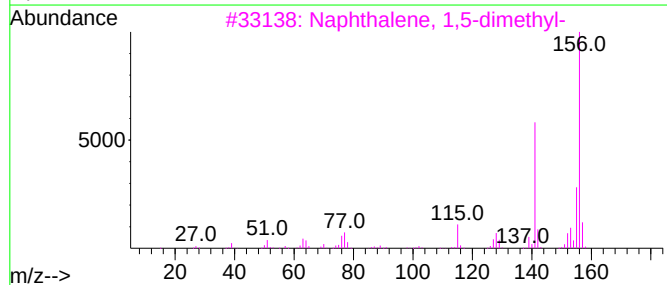
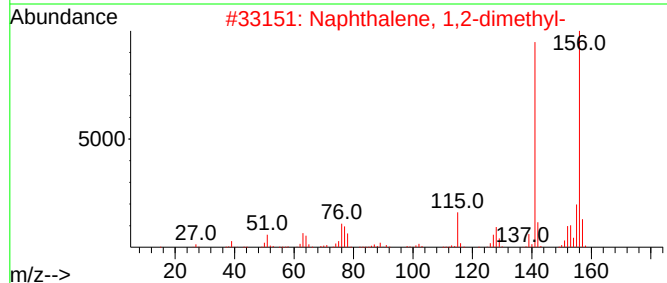
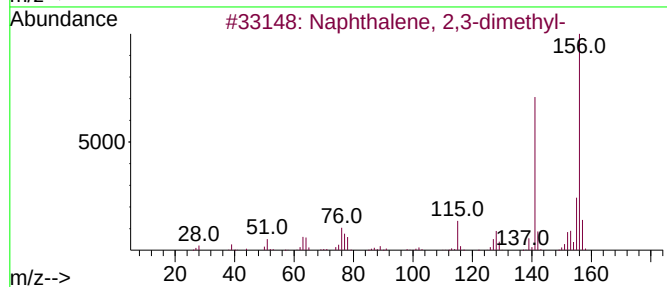
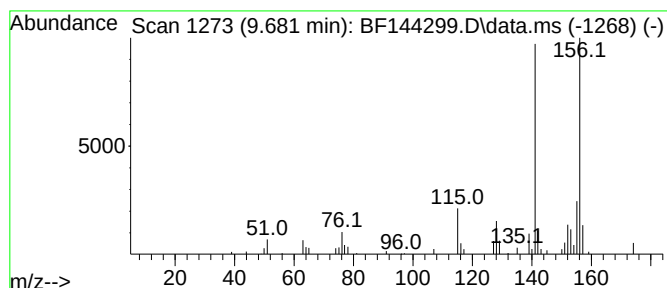
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Naphthalene, 2,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.681	3.08 ng	49180	Acenaphthene-d10	9.904

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2		Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	96
3		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	96
4		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
5		Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112025\
Data File : BF144299.D
Acq On : 20 Nov 2025 17:27
Operator : RC/JU
Sample : Q3681-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

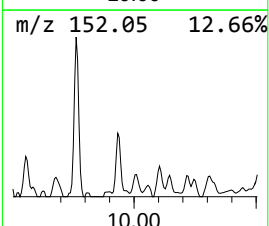
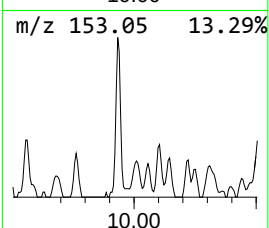
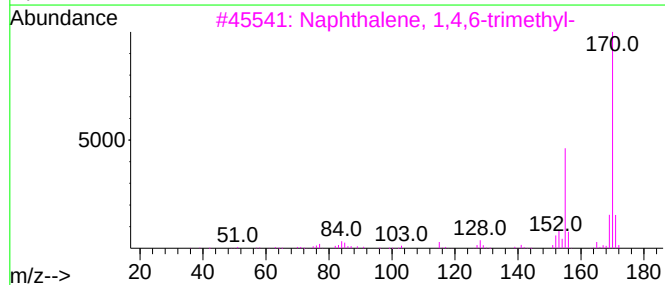
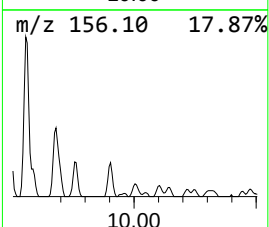
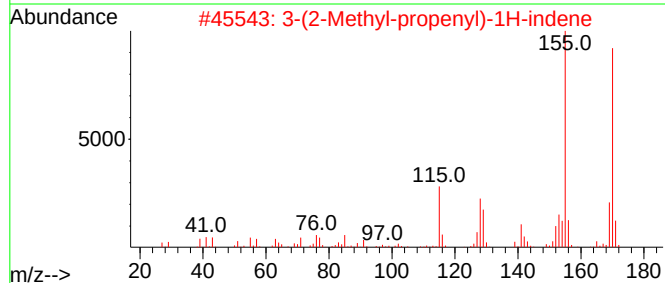
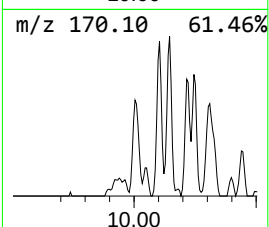
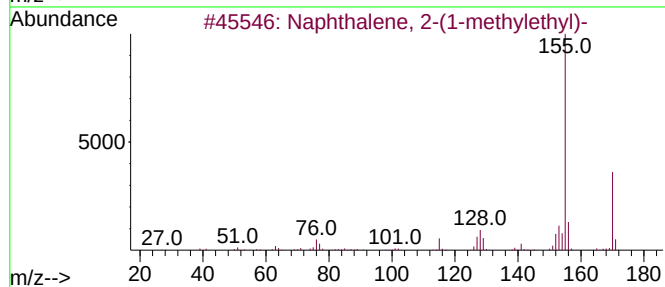
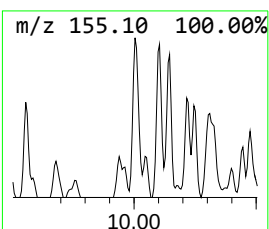
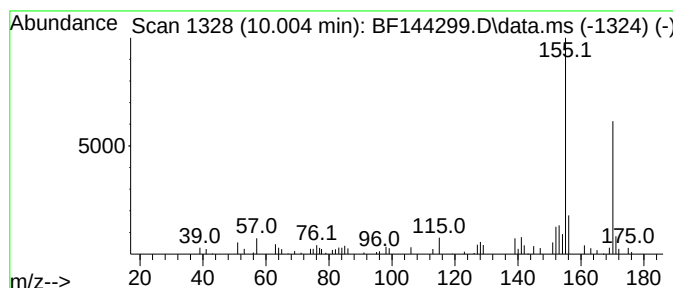
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Naphthalene, 2-(1-methyleth... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD			R.T.
10.004	2.17 ng	34617	Acenaphthene-d10			9.904
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Naphthalene, 2-(1-methylethyl)-		170	C13H14	002027-17-0	91
2	3-(2-Methyl-propenyl)-1H-indene		170	C13H14	1000187-78-5	87
3	Naphthalene, 1,4,6-trimethyl-		170	C13H14	002131-42-2	86
4	Naphthalene, 1,4,5-trimethyl-		170	C13H14	002131-41-1	80
5	4,6,8-Trimethylazulene		170	C13H14	000941-81-1	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112025\
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Operator : RC/JU
Sample : Q3681-01
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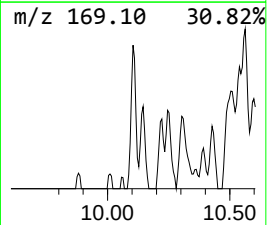
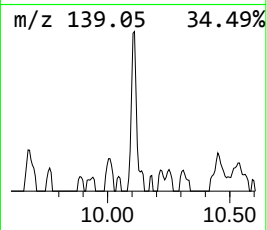
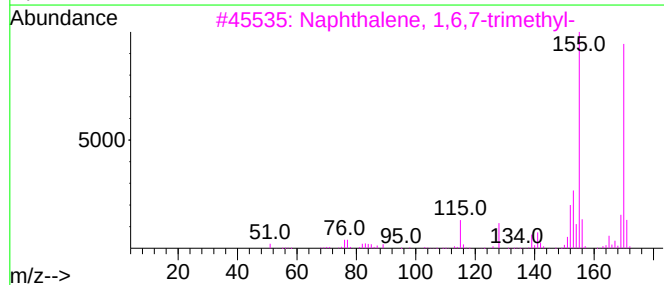
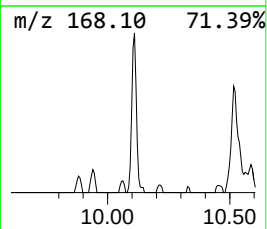
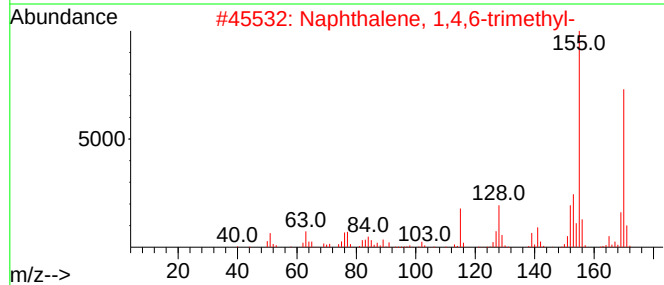
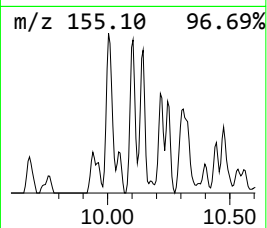
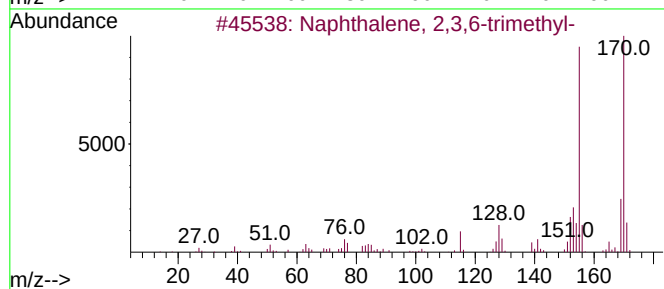
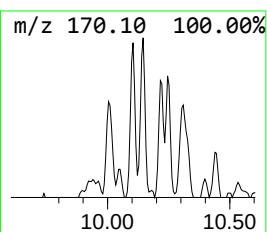
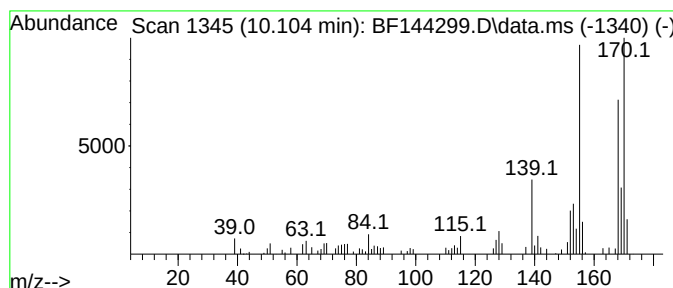
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Naphthalene, 2,3,6-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
10.104	3.81 ng	60749	Acenaphthene-d10		9.904	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3,6-trimethyl-		170	C13H14	000829-26-5	95
2	Naphthalene, 1,4,6-trimethyl-		170	C13H14	002131-42-2	94
3	Naphthalene, 1,6,7-trimethyl-		170	C13H14	002245-38-7	90
4	Naphthalene, 1,4,5-trimethyl-		170	C13H14	002131-41-1	86
5	4,6,8-Trimethylazulene		170	C13H14	000941-81-1	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112025\
Data File : BF144299.D
Acq On : 20 Nov 2025 17:27
Operator : RC/JU
Sample : Q3681-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

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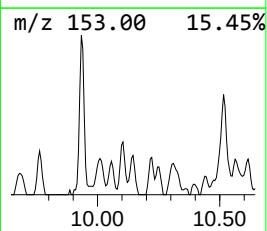
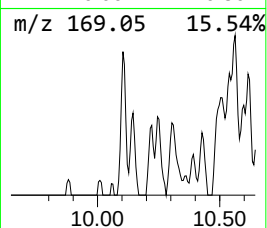
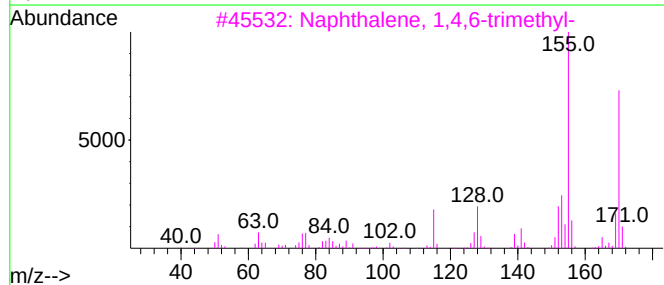
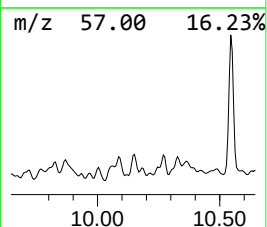
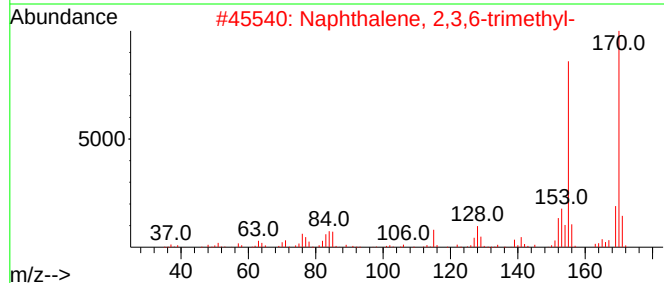
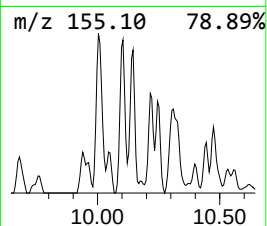
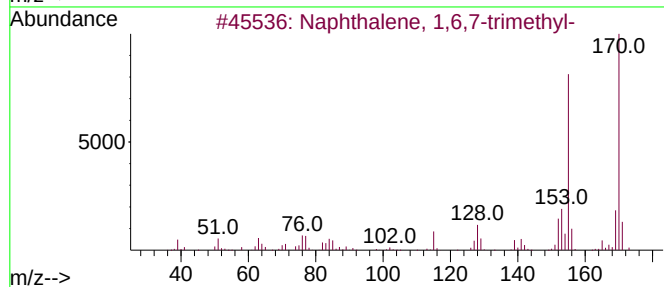
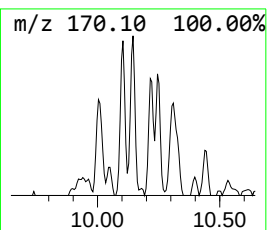
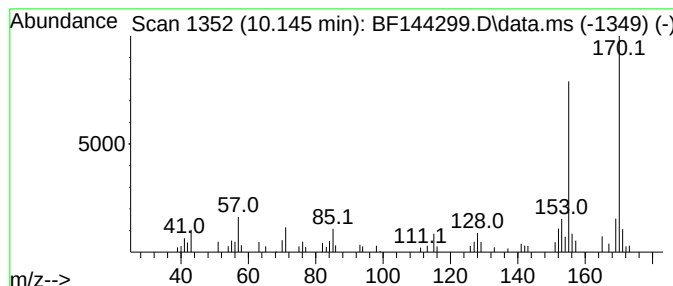
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Naphthalene, 1,6,7-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.145	3.16 ng	50481	Acenaphthene-d10	9.904

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	96
2		Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	96
3		Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	93
4		1-(2,2-Dimethylcyclopropyl)-2-ph...	170	C13H14	1000294-91-1	90
5		4,6,8-Trimethylazulene	170	C13H14	000941-81-1	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112025\
Data File : BF144299.D
Acq On : 20 Nov 2025 17:27
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Misc :
ALS Vial : 11 Sample Multiplier: 1

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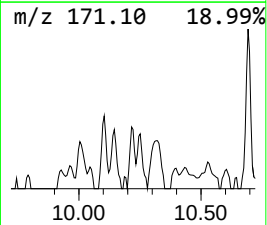
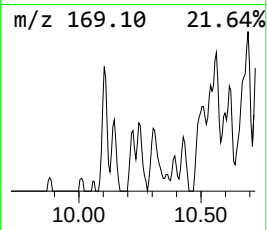
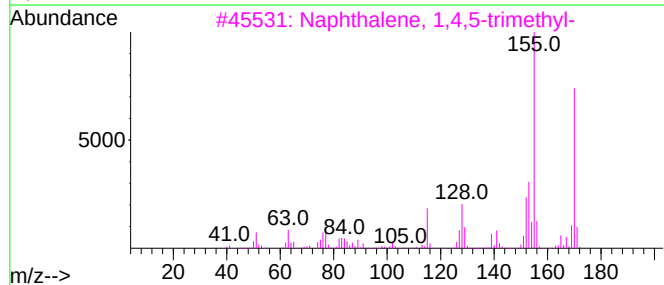
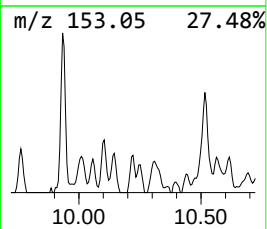
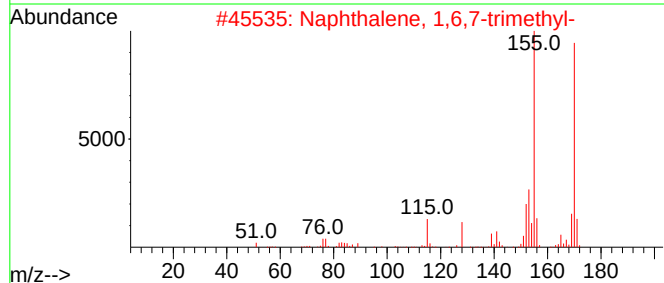
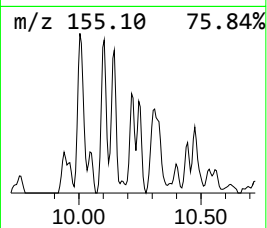
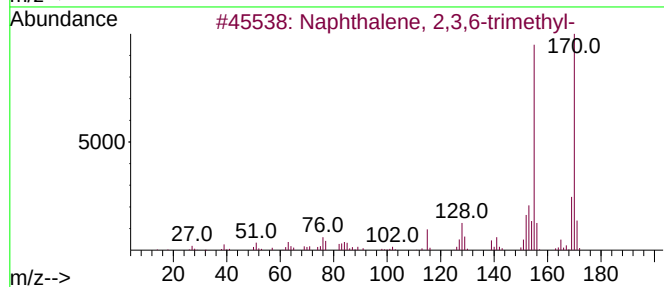
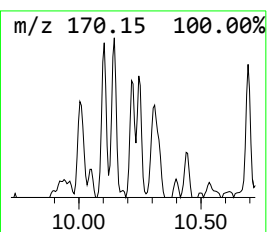
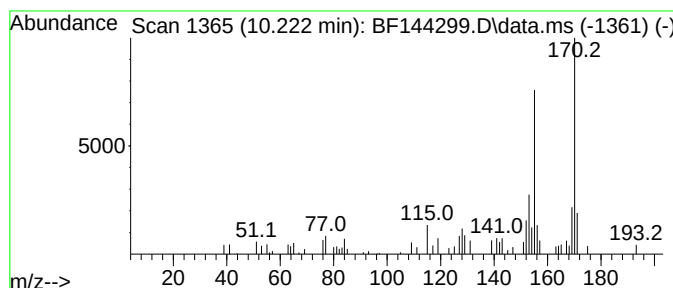
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Naphthalene, 1,4,5-trimethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.222	2.09 ng	33318	Acenaphthene-d10	9.904

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	96
2			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	95
3			Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	95
4			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	94
5			1-(2,2-Dimethylcyclopropyl)-2-ph...	170	C13H14	1000294-91-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112025\
Data File : BF144299.D
Acq On : 20 Nov 2025 17:27
Operator : RC/JU
Sample : Q3681-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
ETGI-354

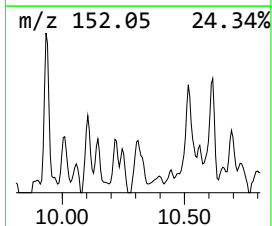
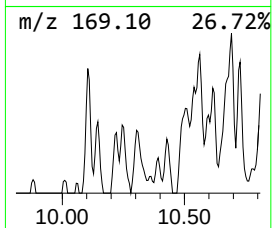
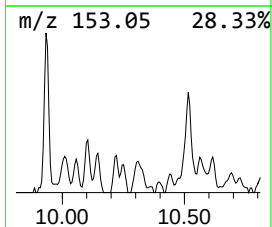
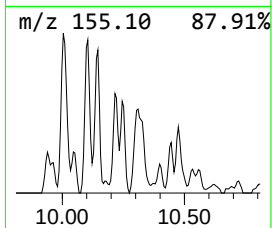
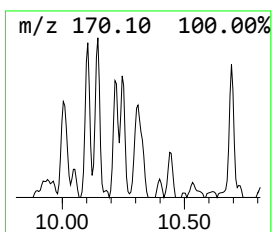
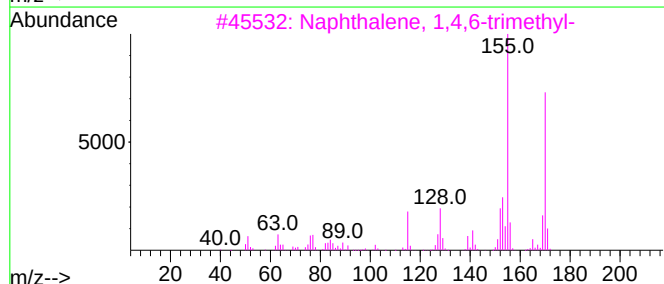
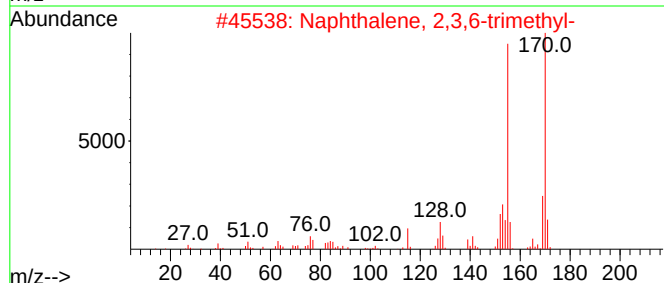
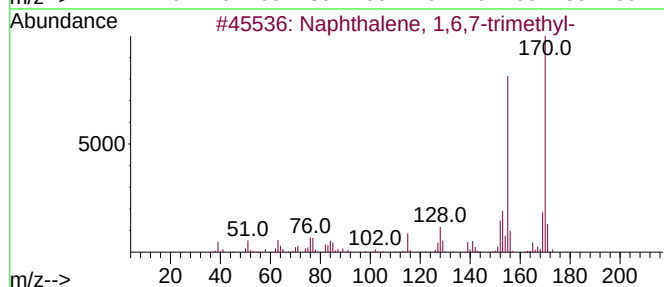
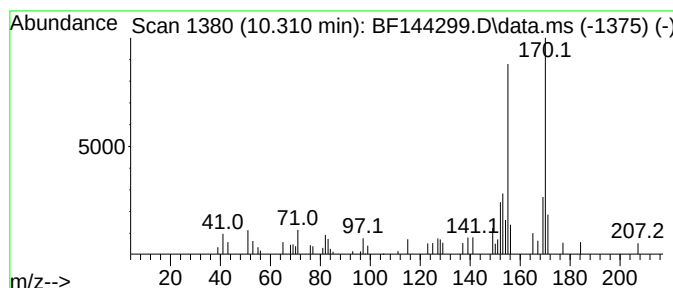
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Naphthalene, 1,4,6-trimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.310	2.67 ng	42640	Acenaphthene-d10	9.904

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	93
2		Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	93
3		Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	90
4		4,6,8-Trimethylazulene	170	C13H14	000941-81-1	83
5		Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112025\
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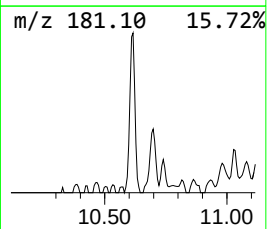
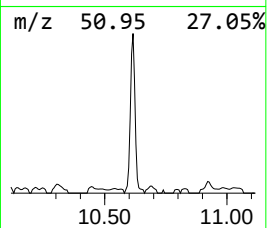
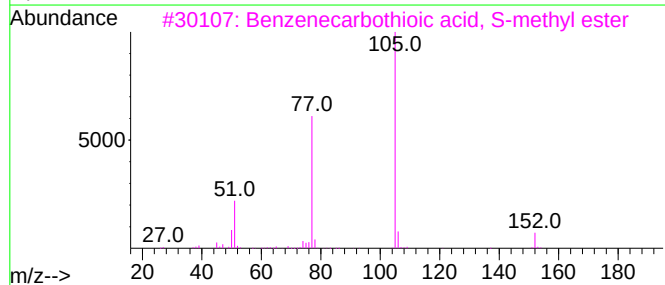
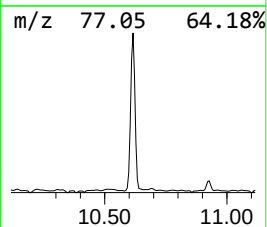
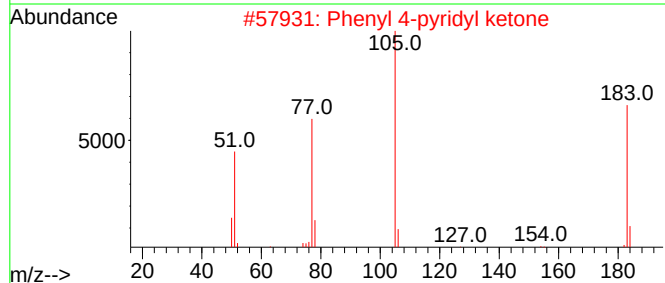
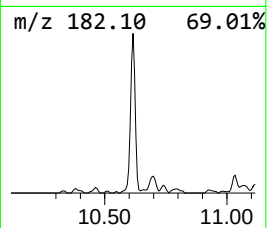
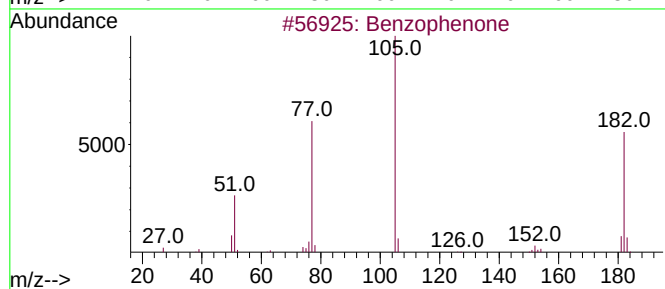
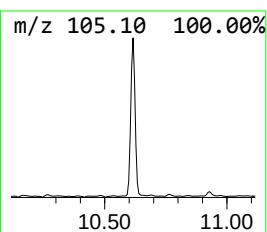
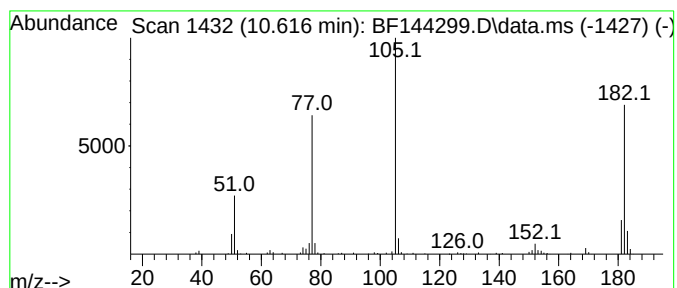
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.616	8.50 ng	135770	Acenaphthene-d10	9.904	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	97
2	Phenyl 4-pyridyl ketone	183	C12H9NO	014548-46-0	58
3	Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	53
4	Benzeneacetic acid, .alpha.-oxo-...	164	C9H8O3	015206-55-0	43
5	Benzenebutanoic acid, .gamma.-oxo-	178	C10H10O3	002051-95-8	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112025\
Data File : BF144299.D
Acq On : 20 Nov 2025 17:27
Operator : RC/JU
Sample : Q3681-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ETGI-354

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110525.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
Phenol, 2,4-dib...	9.157	2.8	ng	43973	3	9.904	319272	20.0
Naphthalene, 1,...	9.486	4.2	ng	67286	3	9.904	319272	20.0
Naphthalene, 2,...	9.557	6.1	ng	97287	3	9.904	319272	20.0
Tetratetracontane	9.616	2.7	ng	43420	3	9.904	319272	20.0
Naphthalene, 2,...	9.681	3.1	ng	49180	3	9.904	319272	20.0
Naphthalene, 2-...	10.004	2.2	ng	34617	3	9.904	319272	20.0
Naphthalene, 2,...	10.104	3.8	ng	60749	3	9.904	319272	20.0
Naphthalene, 1,...	10.145	3.2	ng	50481	3	9.904	319272	20.0
Naphthalene, 1,...	10.222	2.1	ng	33318	3	9.904	319272	20.0
Naphthalene, 1,...	10.310	2.7	ng	42640	3	9.904	319272	20.0
Benzophenone	10.616	8.5	ng	135770	3	9.904	319272	20.0