

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF082020\
 Data File : BF121337.D
 Acq On : 20 Aug 2020 10:08
 Operator : JU/CG
 Sample : L3712-11 5X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CHRT-25433

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 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

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081920.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.293	541	545	558	rBV	450139	503947	29.43%	2.977%
2	6.328	718	721	729	rBV	498056	524694	30.64%	3.100%
3	6.687	779	782	786	rBV	1027318	873583	51.01%	5.161%
4	7.251	875	878	883	rBV	409508	323670	18.90%	1.912%
5	7.975	997	1001	1004	rBV	1403897	1177438	68.75%	6.957%
6	8.934	1160	1164	1166	rBV5	22498	31118	1.82%	0.184%
7	8.998	1171	1175	1178	rBV3	47317	50077	2.92%	0.296%
8	9.045	1179	1183	1187	rBV	824356	654692	38.23%	3.868%
9	9.122	1193	1196	1202	rBV5	65746	105742	6.17%	0.625%
10	9.357	1233	1236	1238	rBV3	59323	70003	4.09%	0.414%
11	9.392	1239	1242	1246	rVB	139077	126232	7.37%	0.746%
12	9.434	1246	1249	1250	rBV	156960	151898	8.87%	0.897%
13	9.586	1273	1275	1278	rBV2	133325	174061	10.16%	1.028%
14	9.634	1280	1283	1286	rVB	307487	275067	16.06%	1.625%
15	9.669	1286	1289	1292	rBV3	90974	128031	7.48%	0.756%
16	9.722	1292	1298	1302	rBV	1635870	1712597	100.00%	10.118%
17	9.986	1341	1343	1346	rBV3	157603	169404	9.89%	1.001%
18	10.092	1359	1361	1364	rBV4	143185	151198	8.83%	0.893%
19	10.128	1365	1367	1370	rBV2	318740	318599	18.60%	1.882%
20	10.386	1408	1411	1416	rBV	339466	356536	20.82%	2.107%
21	10.510	1429	1432	1435	rBV2	402704	454504	26.54%	2.685%
22	10.651	1453	1456	1460	rVB	896050	843884	49.28%	4.986%
23	11.116	1533	1535	1539	rVB	815992	828672	48.39%	4.896%
24	11.210	1548	1551	1554	rBV	992135	933870	54.53%	5.518%
25	12.792	1817	1820	1828	rVB	903824	910780	53.18%	5.381%
26	13.710	1967	1976	1982	rVB3	400774	796561	46.51%	4.706%
27	13.839	1994	1998	2001	rBV	1643903	1521076	88.82%	8.987%
28	14.316	2076	2079	2080	rBV	46928	39919	2.33%	0.236%
29	14.339	2080	2083	2092	rVB2	104384	135737	7.93%	0.802%
30	14.604	2124	2128	2132	rBV2	97207	131531	7.68%	0.777%
31	14.674	2138	2140	2146	rVB	34327	40991	2.39%	0.242%
32	14.845	2166	2169	2178	rVB	37666	65856	3.85%	0.389%
33	14.921	2178	2182	2186	rBV	185194	201622	11.77%	1.191%
34	15.245	2232	2237	2241	rBV	1265122	1485946	86.77%	8.779%

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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

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35	15.274	2241	2242	2248	rVB	106494	89970	5.25%	0.532%
36	15.674	2305	2310	2315	rBV	118620	167733	9.79%	0.991%
37	15.733	2317	2320	2326	rVV6	31552	63691	3.72%	0.376%
38	15.815	2330	2334	2341	rVB6	36217	62963	3.68%	0.372%
39	16.133	2383	2388	2399	rVB2	63959	114677	6.70%	0.678%
40	16.668	2474	2479	2486	rVB3	43067	86609	5.06%	0.512%
41	17.115	2553	2555	2569	rVB10	26210	70304	4.11%	0.415%

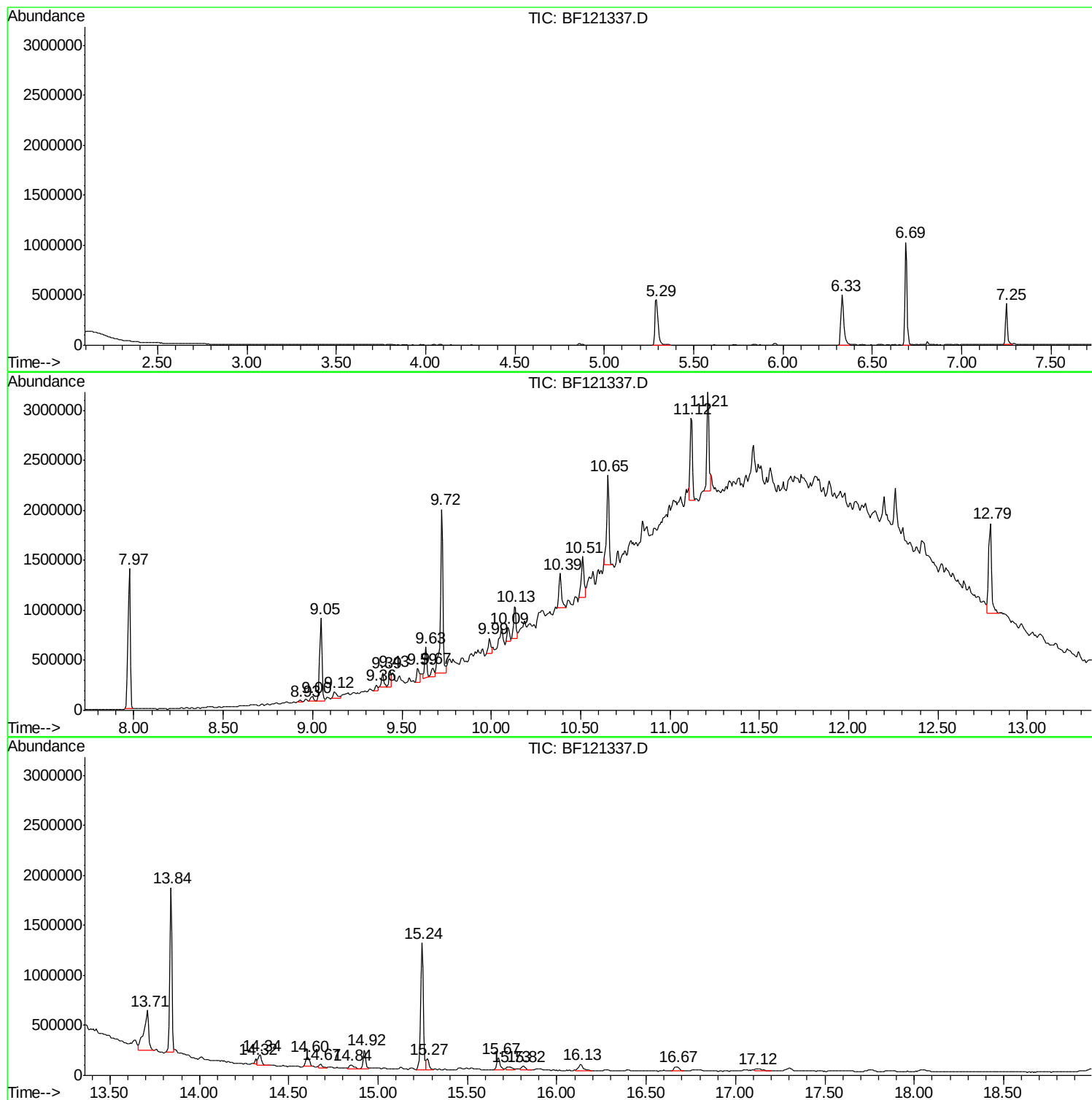
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF081920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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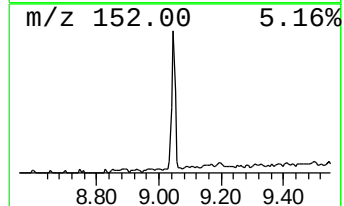
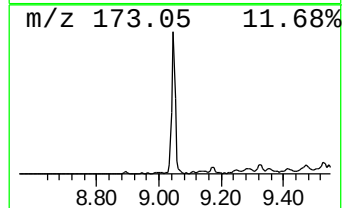
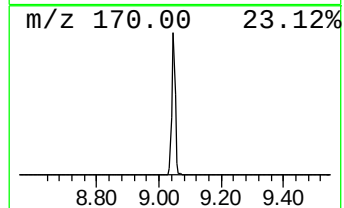
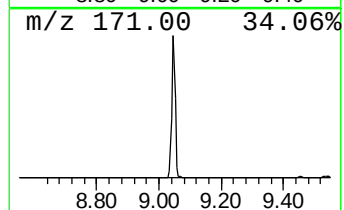
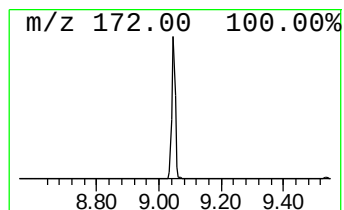
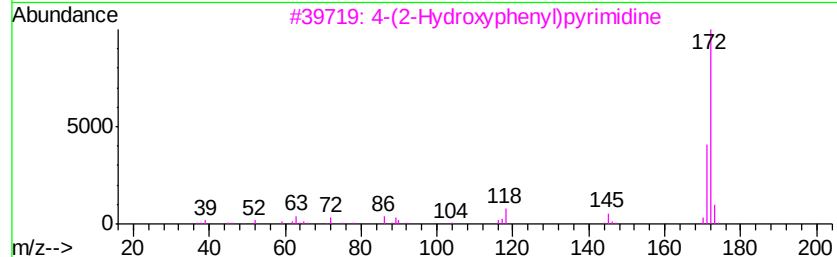
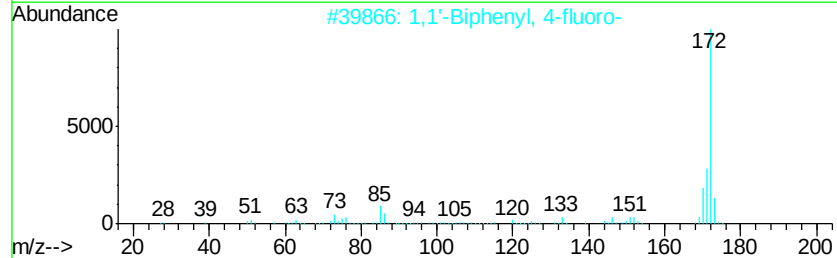
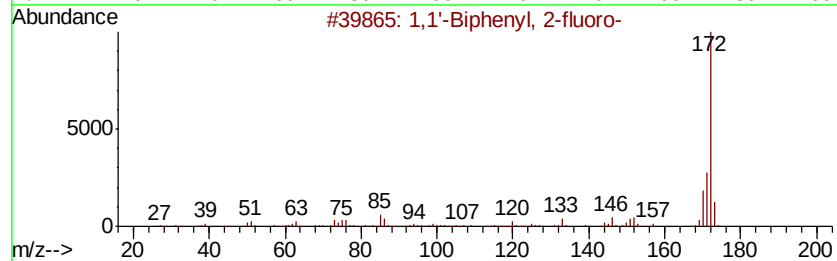
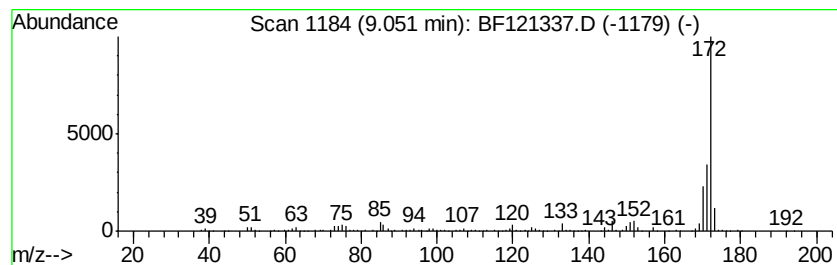
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF081920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 1,1'-Biphenyl, 2-fluoro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.		
9.05	7.65 ng	654692	Acenaphthene-d10		9.72		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'	-Biphenyl,	2-fluoro-	172	C12H9F	000321-60-8	96
2	1,1'	-Biphenyl,	4-fluoro-	172	C12H9F	000324-74-3	93
3	4-(2-Hydroxyphenyl)	pyrimidine		172	C10H8N2O	068535-55-7	45
4	4-Pyrimidinol,	5-phenyl-		172	C10H8N2O	022433-69-8	42
5	2-(2-Hydroxyphenyl)	pyrimidine		172	C10H8N2O	064435-20-7	38



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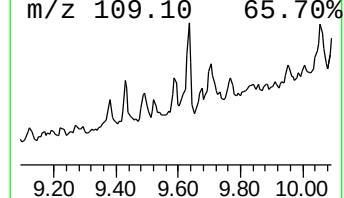
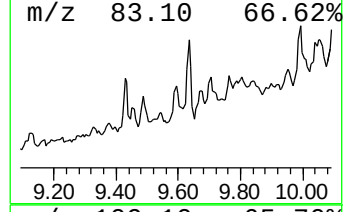
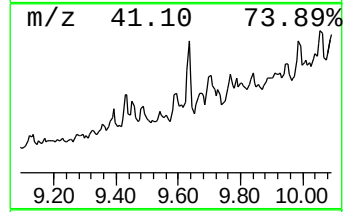
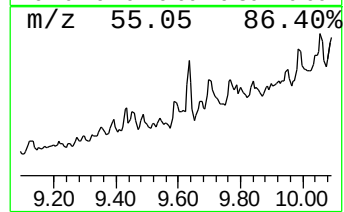
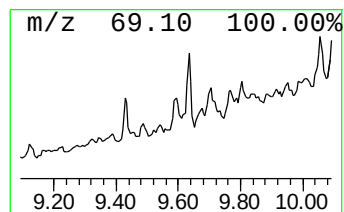
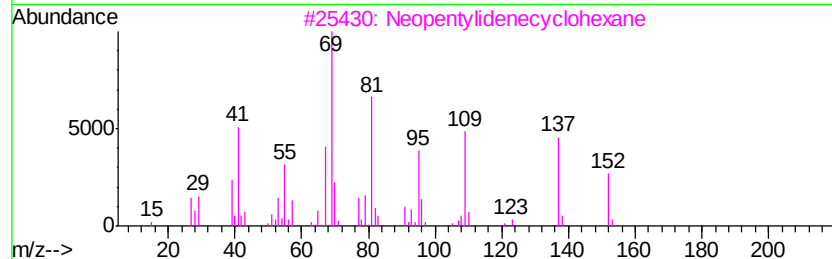
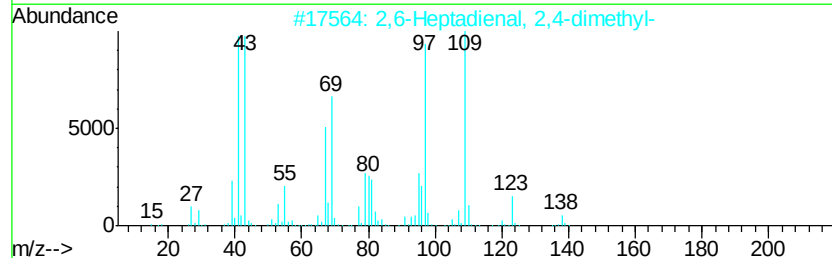
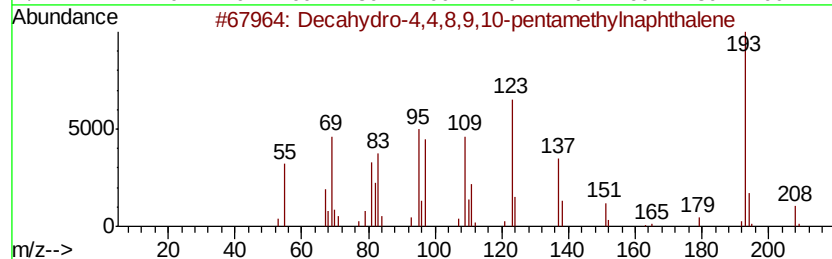
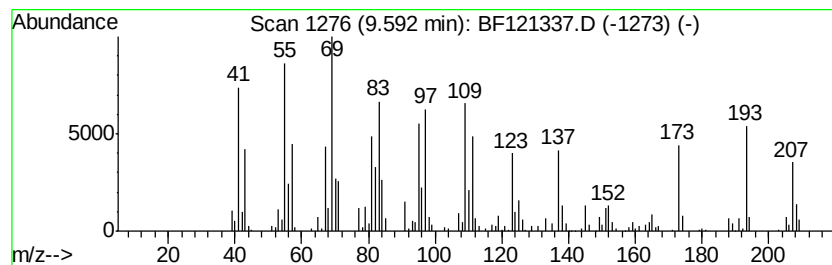
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TIC Integration Parameters: LSCINT.P

Peak Number 2 Decahydro-4,4,8,9,10-pentam... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.		
9.59	2.03 ng	174061	Acenaphthene-d10	9.72		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	90
2		2,6-Heptadienal, 2,4-dimethyl-	138	C9H14O	085136-08-9	42
3		Neopentylidenecyclohexane	152	C11H20	039546-80-0	25
4		2-Anthracenamine	193	C14H11N	000613-13-8	25
5		1H-Pyrazole, 1-methyl-4-(4,4,5,5...	208	C10H17BN2O2	1000319-69-0	25



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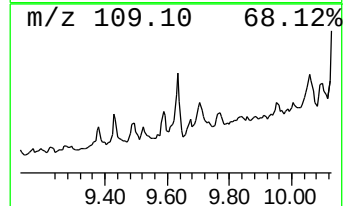
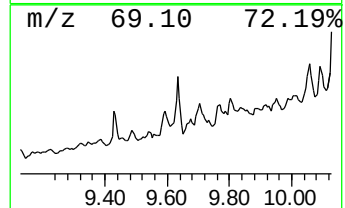
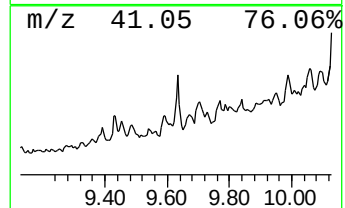
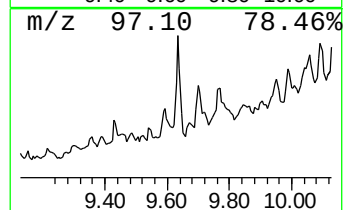
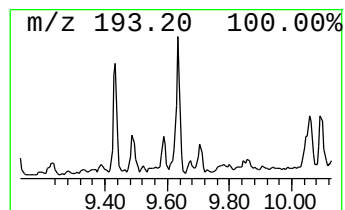
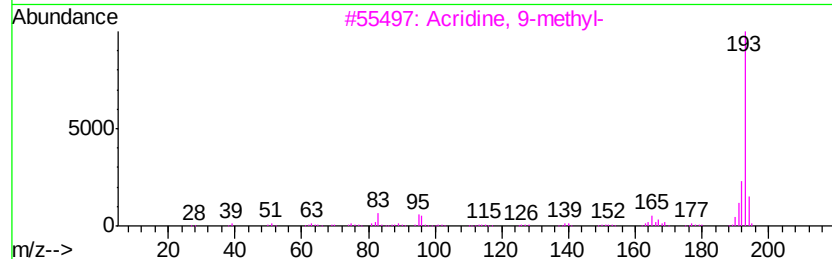
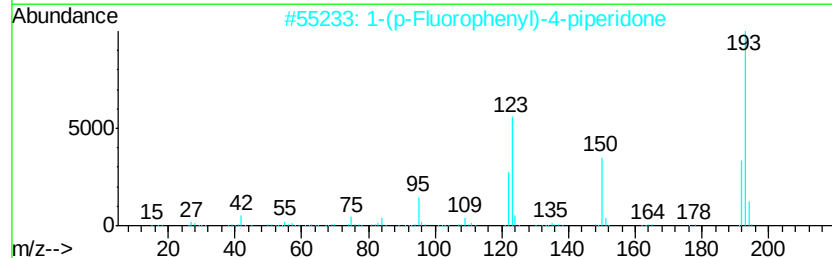
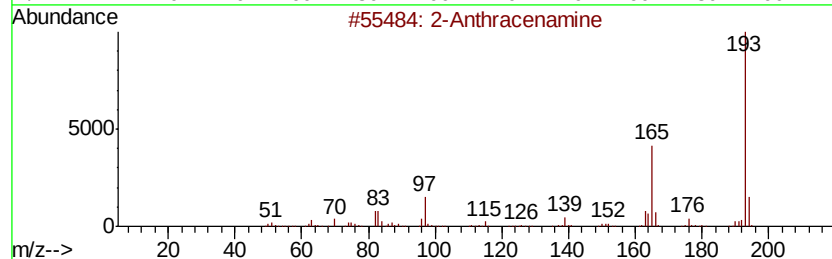
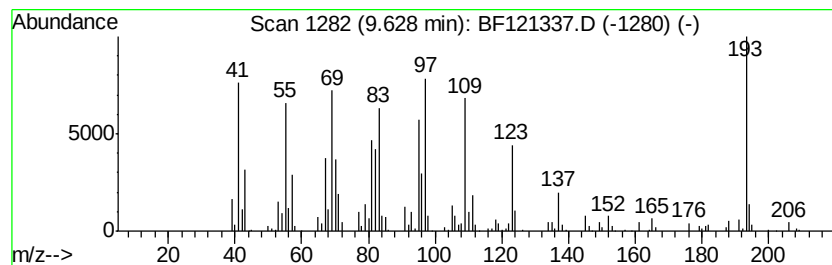
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown9.63 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.63	3.21 ng	275067	Acenaphthene-d10	9.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Anthracenamine	193	C14H11N	000613-13-8	42
2		1-(p-Fluorophenyl)-4-piperidone	193	C11H12FNO	1000238-56-7	27
3		Acridine, 9-methyl-	193	C14H11N	000611-64-3	22
4		Bicyclo[3.1.1]heptane, 2,6,6-tri...	138	C10H18	000473-55-2	20
5		Cyclohexane, 1-(cyclohexylmethyl...	194	C14H26	054823-97-1	18



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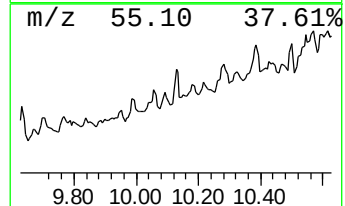
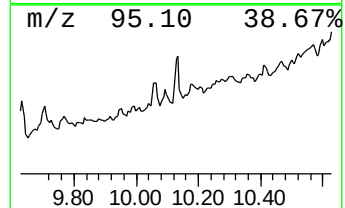
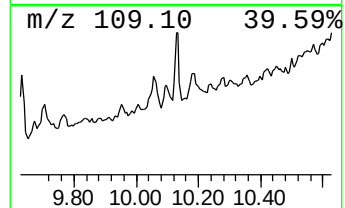
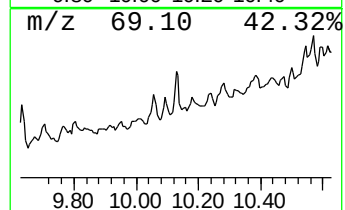
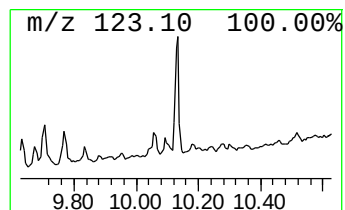
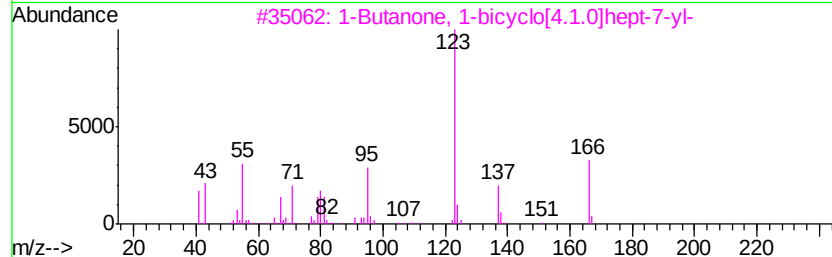
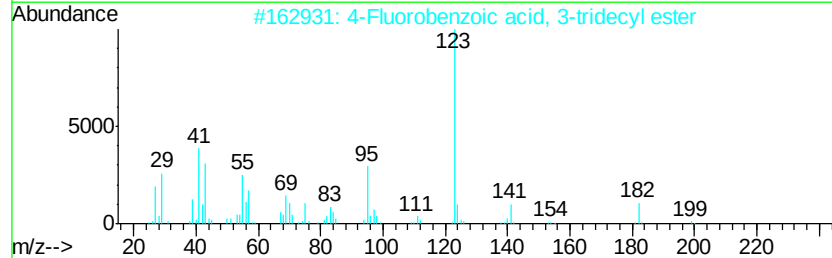
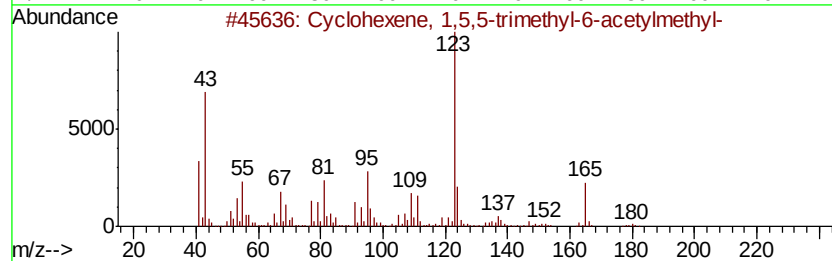
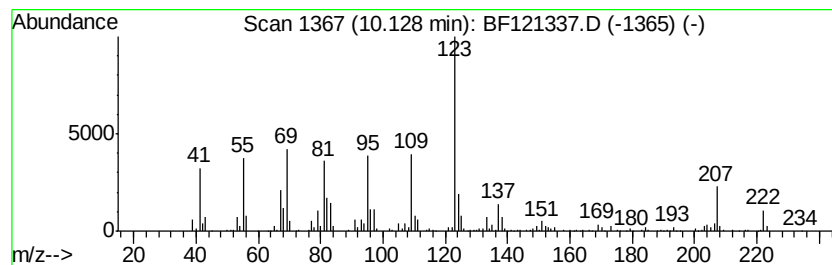
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Cyclohexene, 1,5,5-trimethy... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.13	3.72 ng	318599	Acenaphthene-d10	9.72		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexene, 1,5,5-trimethyl-6-a...	180	C12H20O	211563-96-1	53
2		4-Fluorobenzoic acid, 3-tridecyl...	322	C20H31FO2	1000280-67-9	47
3		1-Butanone, 1-bicyclo[4.1.0]hept...	166	C11H18O	054764-62-4	47
4		3-Fluorobenzoic acid, 2-tridecyl...	322	C20H31FO2	1000280-60-2	47
5		2-Fluorobenzoic acid, 2-ethylcyc...	250	C15H19FO2	1000278-98-1	43



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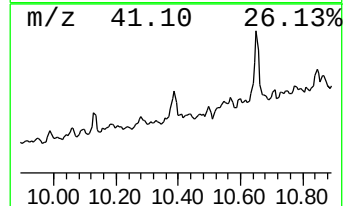
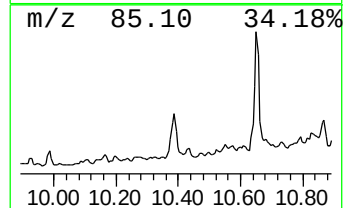
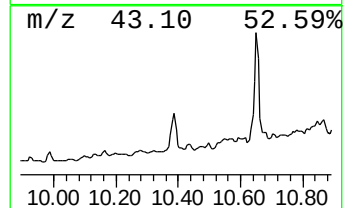
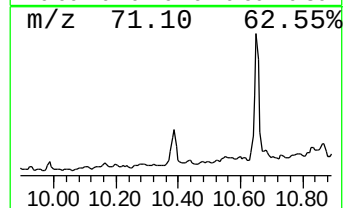
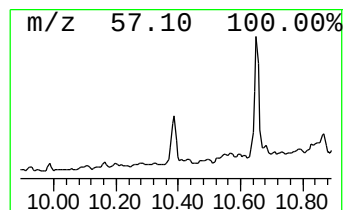
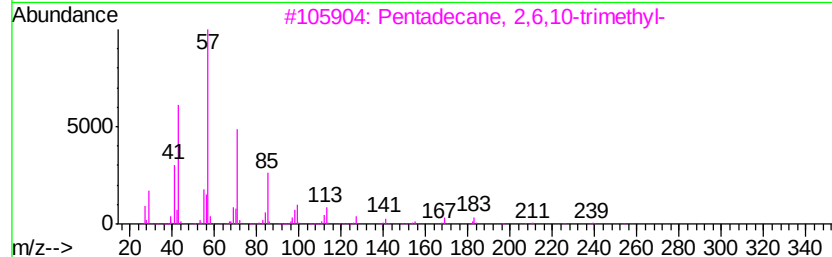
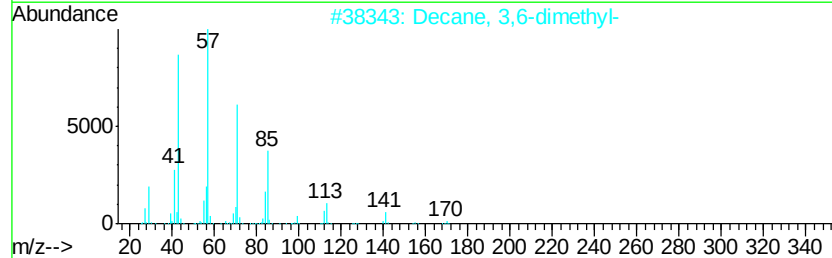
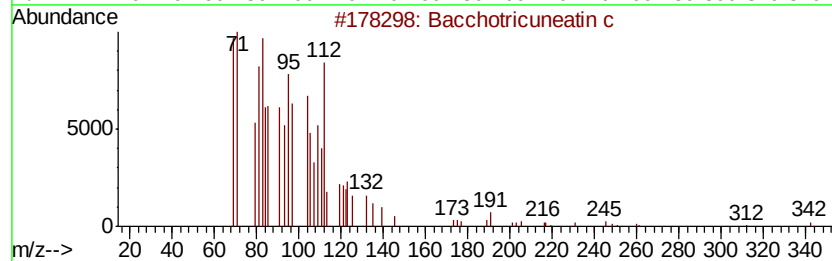
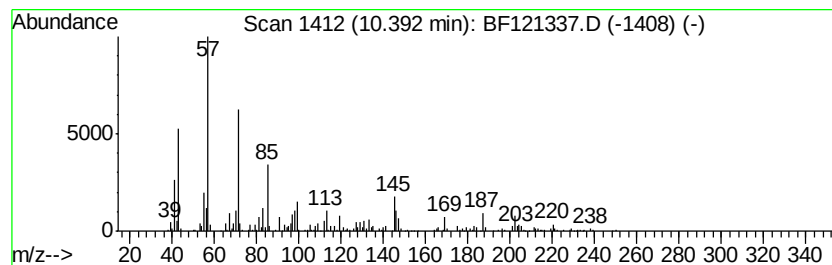
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Bacchotricuneatin c Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.39	4.16 ng	356536	Acenaphthene-d10	9.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bacchotricuneatin c	342	C20H22O5	066563-30-2	96
2		Decane, 3,6-dimethyl-	170	C12H26	017312-53-7	81
3		Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	80
4		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	68
5		Tridecane	184	C13H28	000629-50-5	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082020\
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 Acq On : 20 Aug 2020 10:08
 Operator : JU/CG
 Sample : L3712-11 5X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

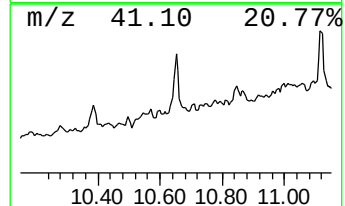
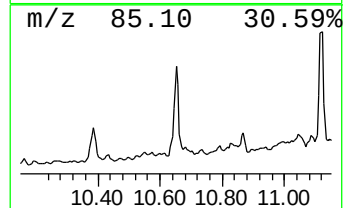
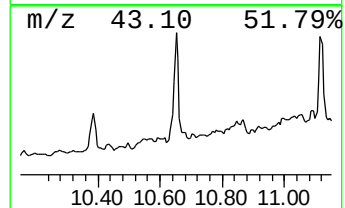
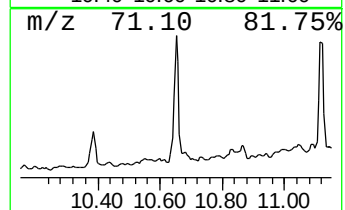
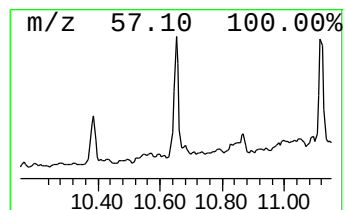
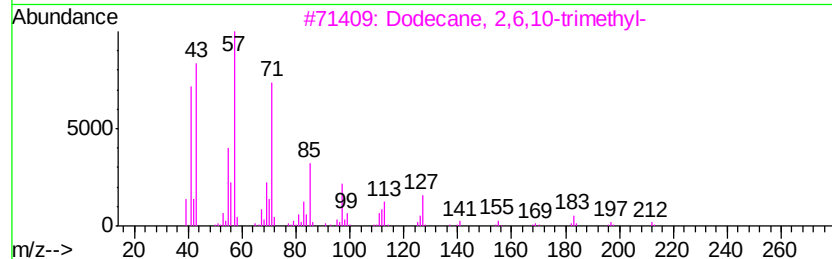
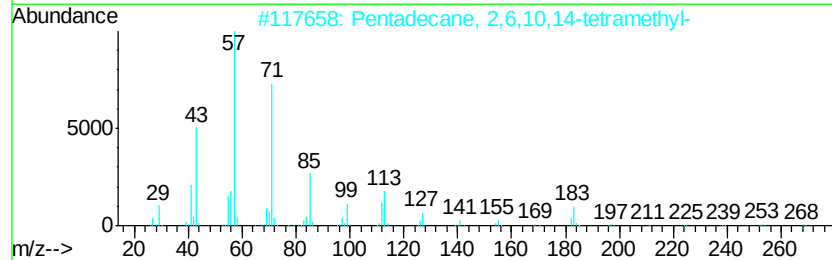
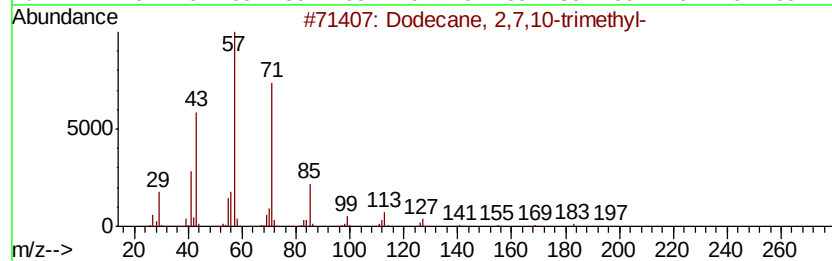
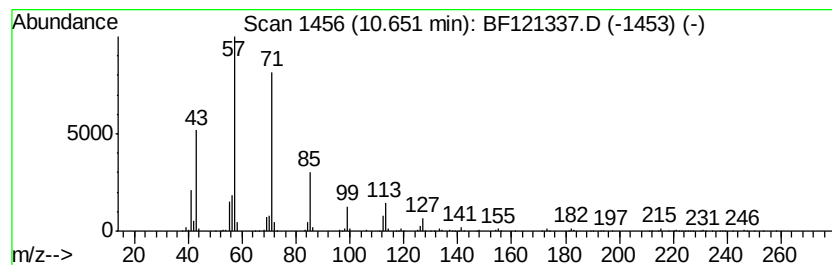
Instrument :
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 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF081920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 6 Dodecane, 2,7,10-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.65	18.07 ng	843884	Phenanthrene-d10	11.21	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	86
2	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	83
3	Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	78
4	Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	72
5	Undecane, 4,8-dimethyl-	184	C13H28	017301-33-6	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082020\
Data File : BF121337.D
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ClientSampleId :
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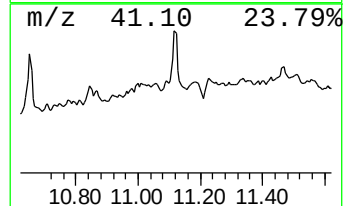
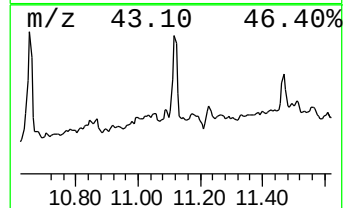
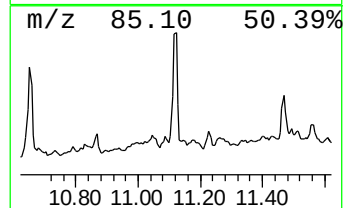
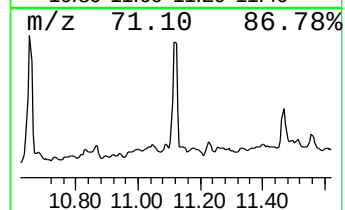
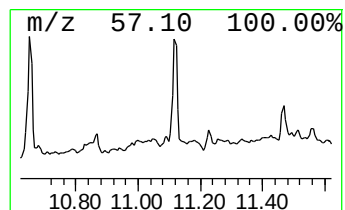
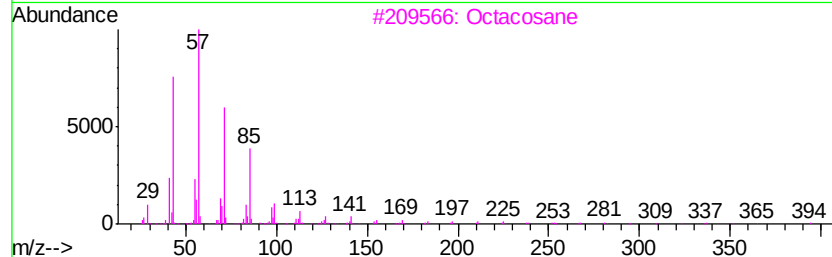
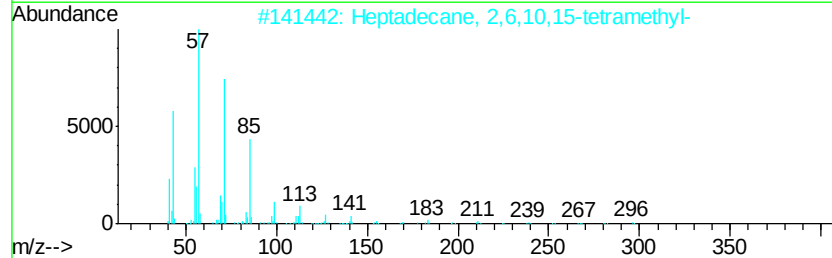
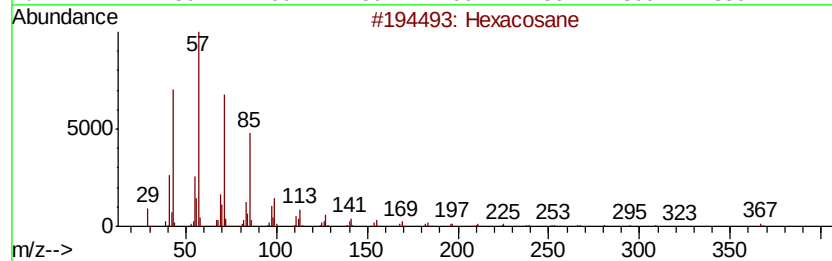
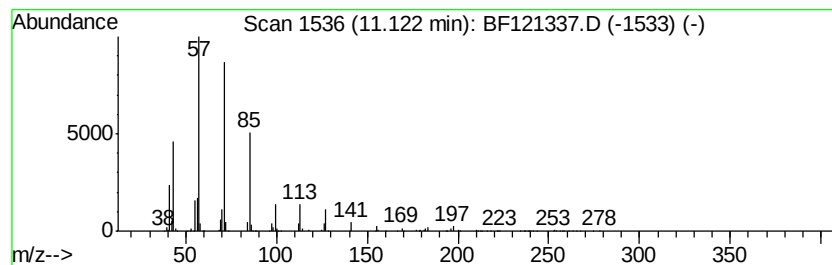
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Hexacosane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.12	17.75 ng	828672	Phenanthrene-d10	11.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	90
2		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86
3		Octacosane	394	C28H58	000630-02-4	86
4		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	83
5		Tetracosane	338	C24H50	000646-31-1	83



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Sample : L3712-11 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

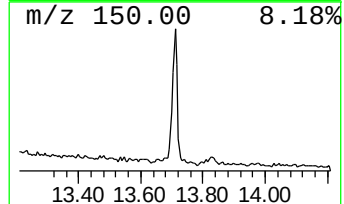
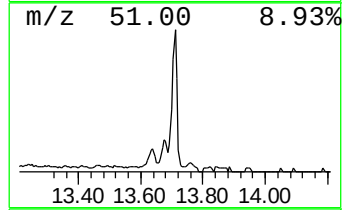
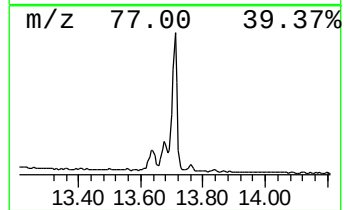
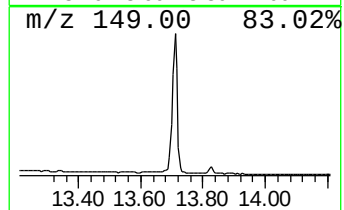
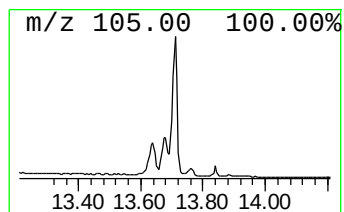
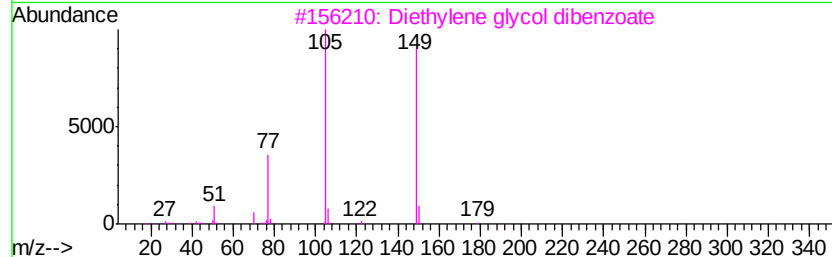
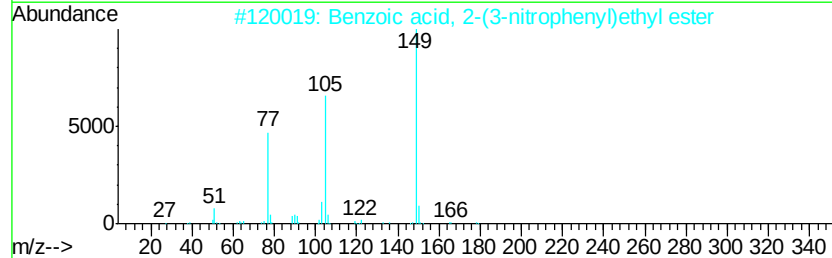
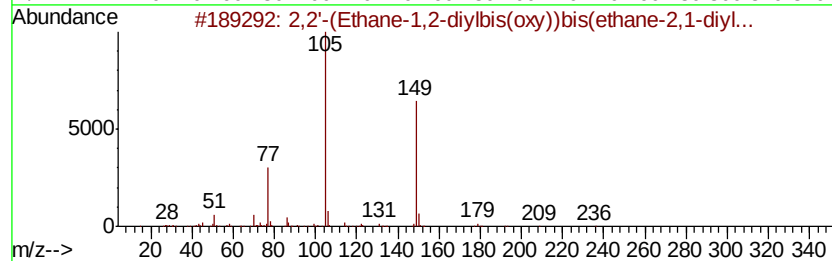
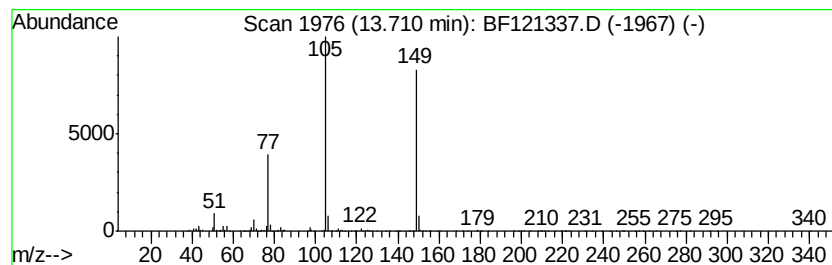
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF081920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 2,2'-(Ethane-1,2-diylbis(ox... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.71	10.47 ng	796561	Chrysene-d12	13.84	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2'-(Ethane-1,2-diylbis(oxv))bi...	358	C20H22O6	1000366-99-4	78
2	Benzoic acid, 2-(3-nitrophenyl)e...	271	C15H13NO4	1000368-22-4	72
3	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	59
4	Phthalic acid, propyl 6-ethyl-3-...	348	C21H32O4	1000315-17-3	43
5	Phthalic acid, dec-2-yl propyl e...	348	C21H32O4	1000356-93-7	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082020\
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Sample : L3712-11 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

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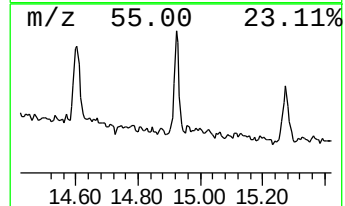
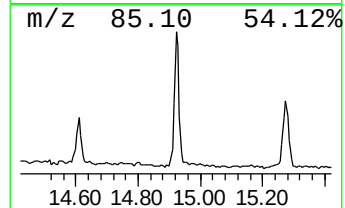
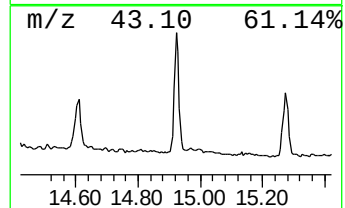
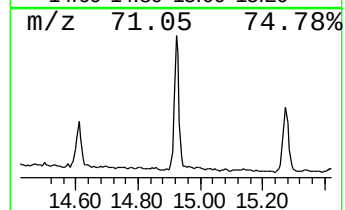
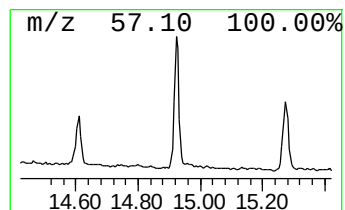
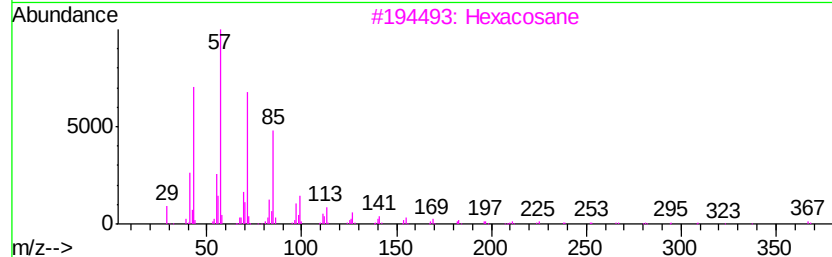
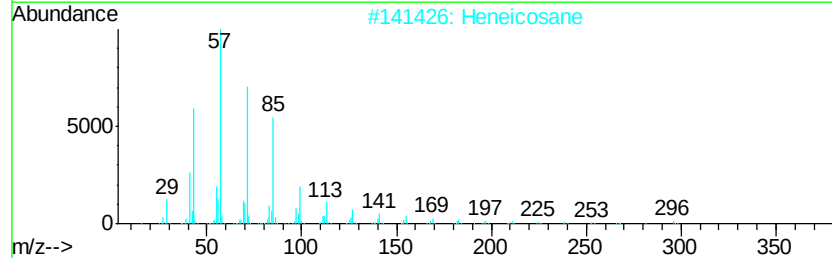
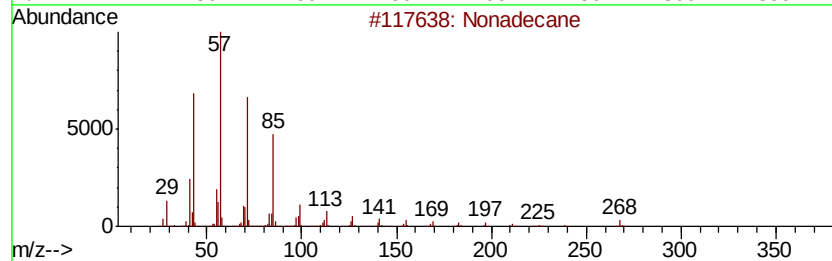
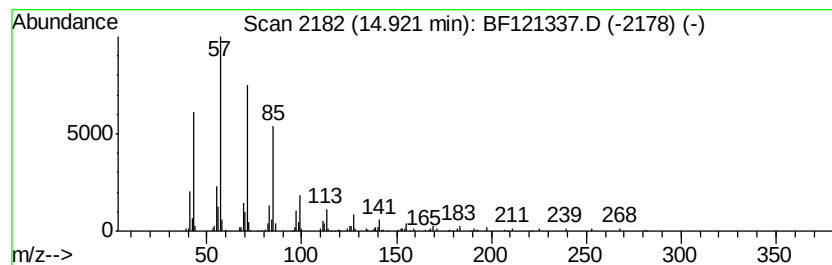
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Nonadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.92	2.71 ng	201622	Perylene-d12	15.24

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	95
2	Heneicosane		296	C21H44	000629-94-7	91
3	Hexacosane		366	C26H54	000630-01-3	91
4	Octacosane		394	C28H58	000630-02-4	91
5	Heptadecane		240	C17H36	000629-78-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082020\
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Misc :
ALS Vial : 4 Sample Multiplier: 1

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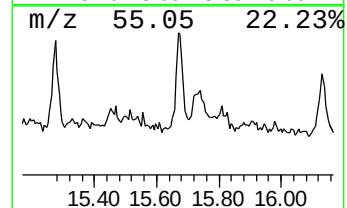
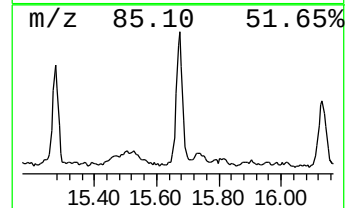
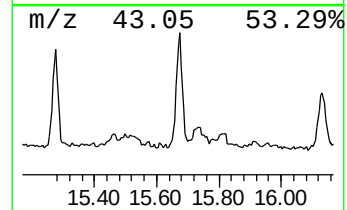
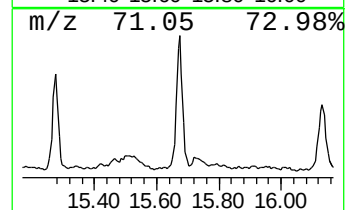
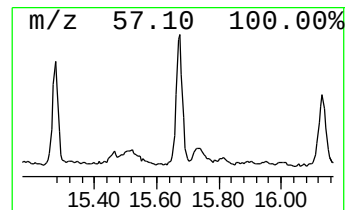
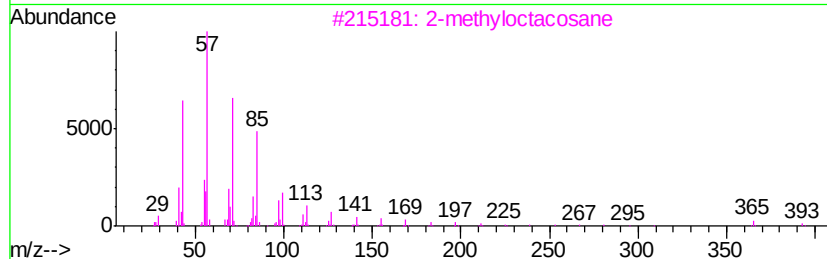
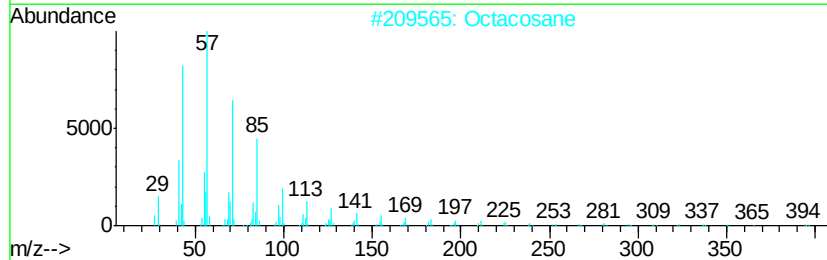
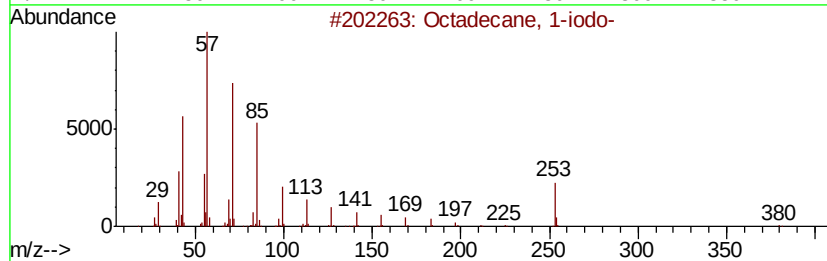
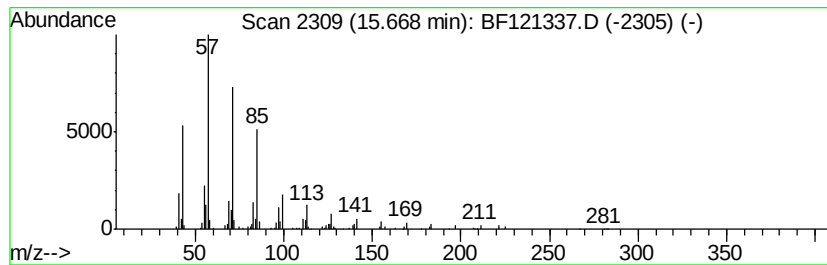
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Octadecane, 1-iodo- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.67	2.26 ng	167733	Perylene-d12	15.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane, 1-iodo-	380	C18H37I	000629-93-6	91
2		Octacosane	394	C28H58	000630-02-4	91
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF082020\
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1,1'-Biphenyl, 2-...	9.05	7.7	ng	654692	3	9.72	1712600	20.0
Decahydro-4,4,8,9...	9.59	2.0	ng	174061	3	9.72	1712600	20.0
unknown9.63	9.63	3.2	ng	275067	3	9.72	1712600	20.0
Cyclohexene, 1,5,...	10.13	3.7	ng	318599	3	9.72	1712600	20.0
Bacchotricuneatin c	10.39	4.2	ng	356536	3	9.72	1712600	20.0
Dodecane, 2,7,10-...	10.65	18.1	ng	843884	4	11.21	933870	20.0
Hexacosane	11.12	17.8	ng	828672	4	11.21	933870	20.0
2,2'-(Ethane-1,2-...	13.71	10.5	ng	796561	5	13.84	1521080	20.0
Nonadecane	14.92	2.7	ng	201622	6	15.24	1485950	20.0
Octadecane, 1-iodo-	15.67	2.3	ng	167733	6	15.24	1485950	20.0