

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051921\
 Data File : BF123995.D
 Acq On : 19 May 2021 17:49
 Operator : JU/CG
 Sample : M2448-01 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HD-01-051821

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-BF051421.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF123995.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.151	5	11	16	rVB	237596	303626	52.38%	4.455%
2	2.257	23	29	37	rVB3	17148	24636	4.25%	0.362%
3	5.104	511	513	518	rVB	25433	25252	4.36%	0.371%
4	5.510	578	582	586	rBV	585293	495361	85.45%	7.269%
5	6.516	750	753	761	rVB	552798	517449	89.26%	7.593%
6	6.728	786	789	796	rVB	13261	13415	2.31%	0.197%
7	6.904	815	819	822	rBV	440353	331985	57.27%	4.872%
8	7.457	910	913	917	rBV	377263	338111	58.33%	4.961%
9	8.186	1033	1037	1040	rBV	570961	437091	75.40%	6.414%
10	8.210	1040	1041	1044	rVB	21366	10670	1.84%	0.157%
11	9.210	1208	1211	1216	rBV	10509	8823	1.52%	0.129%
12	9.263	1216	1220	1223	rBV	730805	579701	100.00%	8.507%
13	9.339	1231	1233	1237	rBV3	8685	10118	1.75%	0.148%
14	9.651	1284	1286	1291	rBV	79684	79046	13.64%	1.160%
15	9.804	1310	1312	1317	rBV4	7056	9432	1.63%	0.138%
16	9.945	1332	1336	1339	rBV	582138	479817	82.77%	7.041%
17	9.975	1339	1341	1345	rVB3	20017	18836	3.25%	0.276%
18	10.145	1366	1370	1374	rBV4	10445	14173	2.44%	0.208%
19	10.316	1396	1399	1403	rVB4	5089	8572	1.48%	0.126%
20	10.357	1403	1406	1407	rBV2	8353	8329	1.44%	0.122%
21	10.374	1407	1409	1412	rVB2	10826	8468	1.46%	0.124%
22	10.486	1426	1428	1432	rBV3	13276	14730	2.54%	0.216%
23	10.592	1441	1446	1449	rBV3	15089	16382	2.83%	0.240%
24	10.727	1465	1469	1473	rBV	397667	334526	57.71%	4.909%
25	10.863	1485	1492	1494	rBV	29765	38537	6.65%	0.565%
26	11.069	1524	1527	1531	rVB5	7067	7246	1.25%	0.106%
27	11.292	1562	1565	1569	rBV5	9092	13445	2.32%	0.197%
28	11.327	1569	1571	1574	rVB	18195	13762	2.37%	0.202%
29	11.380	1576	1580	1583	rBV4	4605	8092	1.40%	0.119%
30	11.427	1585	1588	1591	rVV	475642	420264	72.50%	6.167%
31	11.451	1591	1592	1596	rVB	34900	23166	4.00%	0.340%
32	11.721	1633	1638	1642	rVB3	13439	13098	2.26%	0.192%
33	11.951	1675	1677	1680	rBV2	22461	20261	3.50%	0.297%
34	12.039	1689	1692	1695	rBV5	10750	13550	2.34%	0.199%
35	12.133	1705	1708	1710	rBV3	8387	9721	1.68%	0.143%
36	12.416	1752	1756	1758	rBV4	9652	9146	1.58%	0.134%

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Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

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37	12.480	1764	1767	1770	rBV4	8329	10296	1.78%	0.151%
38	12.563	1779	1781	1785	rBV3	7401	7139	1.23%	0.105%
39	12.639	1791	1794	1798	rBV	60575	53534	9.23%	0.786%
40	12.733	1807	1810	1814	rBV5	8206	9478	1.63%	0.139%
41	12.868	1828	1833	1836	rBV	53944	60877	10.50%	0.893%
42	13.015	1854	1858	1861	rBV	569687	500682	86.37%	7.347%
43	13.204	1886	1890	1891	rVB4	12448	12190	2.10%	0.179%
44	13.263	1899	1900	1904	rVB4	7937	8345	1.44%	0.122%
45	13.298	1904	1906	1910	rBV4	10381	10669	1.84%	0.157%
46	13.392	1920	1922	1926	rBV4	10176	12202	2.10%	0.179%
47	13.427	1926	1928	1930	rBV3	9100	7418	1.28%	0.109%
48	13.468	1932	1935	1936	rBV3	13715	10067	1.74%	0.148%
49	13.498	1936	1940	1943	rBV6	13140	20347	3.51%	0.299%
50	13.704	1972	1975	1978	rBV5	9261	12399	2.14%	0.182%
51	13.874	2002	2004	2006	rBV3	7524	7160	1.24%	0.105%
52	13.910	2007	2010	2015	rVV2	70681	85628	14.77%	1.257%
53	14.068	2033	2037	2040	rBV	372782	392371	67.69%	5.758%
54	14.151	2047	2051	2053	rBV5	12802	16362	2.82%	0.240%
55	14.233	2063	2065	2071	rVB6	18285	29215	5.04%	0.429%
56	14.292	2071	2075	2077	rBV5	16140	18107	3.12%	0.266%
57	14.492	2107	2109	2113	rBV5	13531	12000	2.07%	0.176%
58	14.551	2116	2119	2123	rVV4	38215	47238	8.15%	0.693%
59	14.851	2168	2170	2171	rBV	14485	8706	1.50%	0.128%
60	15.004	2193	2196	2201	rBV6	15489	23516	4.06%	0.345%
61	15.080	2207	2209	2210	rBV2	11706	7009	1.21%	0.103%
62	15.127	2213	2217	2221	rVV4	30490	49517	8.54%	0.727%
63	15.209	2227	2231	2234	rBV5	27501	35608	6.14%	0.523%
64	15.433	2264	2269	2273	rBV5	26072	40422	6.97%	0.593%
65	15.504	2277	2281	2283	rVV4	24082	34918	6.02%	0.512%
66	15.568	2287	2292	2295	rVV	223466	290388	50.09%	4.261%
67	15.598	2295	2297	2304	rVB8	30352	52557	9.07%	0.771%
68	15.721	2316	2318	2323	rBV6	12193	18628	3.21%	0.273%
69	15.809	2331	2333	2336	rBV4	10781	10009	1.73%	0.147%
70	16.056	2372	2375	2379	rVB6	16840	22659	3.91%	0.333%
71	16.115	2381	2385	2393	rBV6	19463	39154	6.75%	0.575%
72	16.251	2406	2408	2411	rVB4	14430	13976	2.41%	0.205%
73	16.304	2415	2417	2424	rVB7	19964	35119	6.06%	0.515%
74	16.562	2459	2461	2462	rBV2	8762	7160	1.24%	0.105%
75	16.580	2462	2464	2468	rVB5	14865	13597	2.35%	0.200%
76	16.756	2488	2494	2499	rBV9	20498	44067	7.60%	0.647%

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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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	17.139	2557	2559	2562	rBV4	6600	8655	1.49%	0.127%
78	17.368	2594	2598	2600	rBV5	9816	12999	2.24%	0.191%
79	17.515	2621	2623	2625	rBV3	8768	7815	1.35%	0.115%
80	17.574	2631	2633	2636	rBV4	5459	7744	1.34%	0.114%
81	18.098	2718	2722	2723	rBV4	6040	7915	1.37%	0.116%

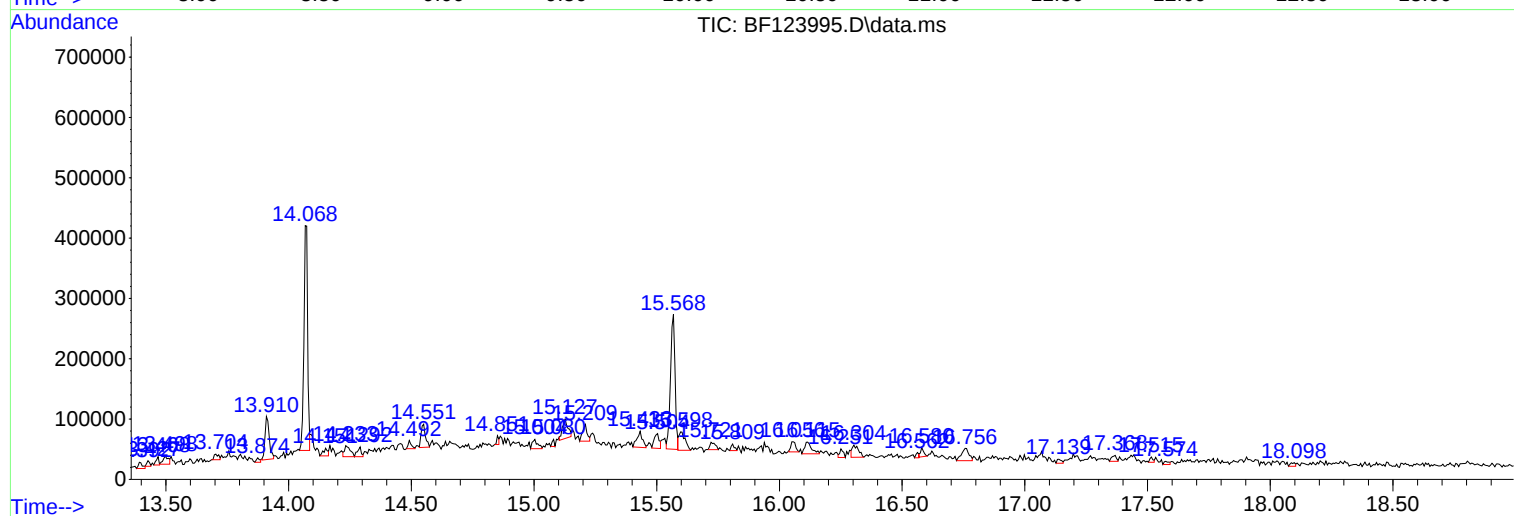
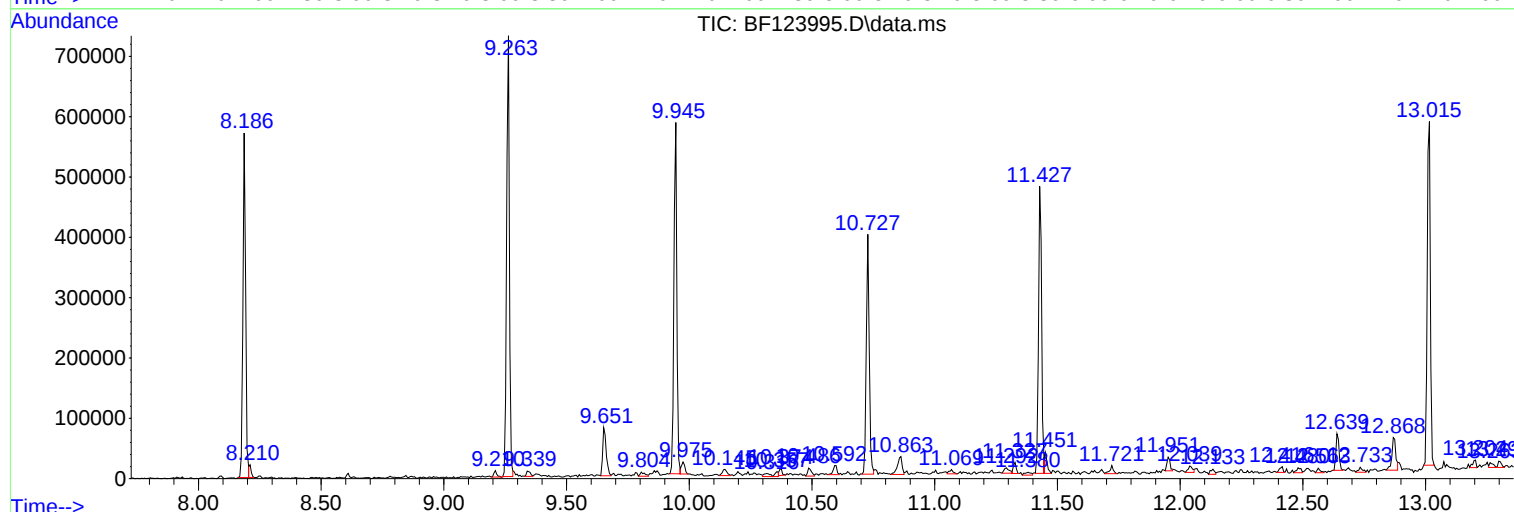
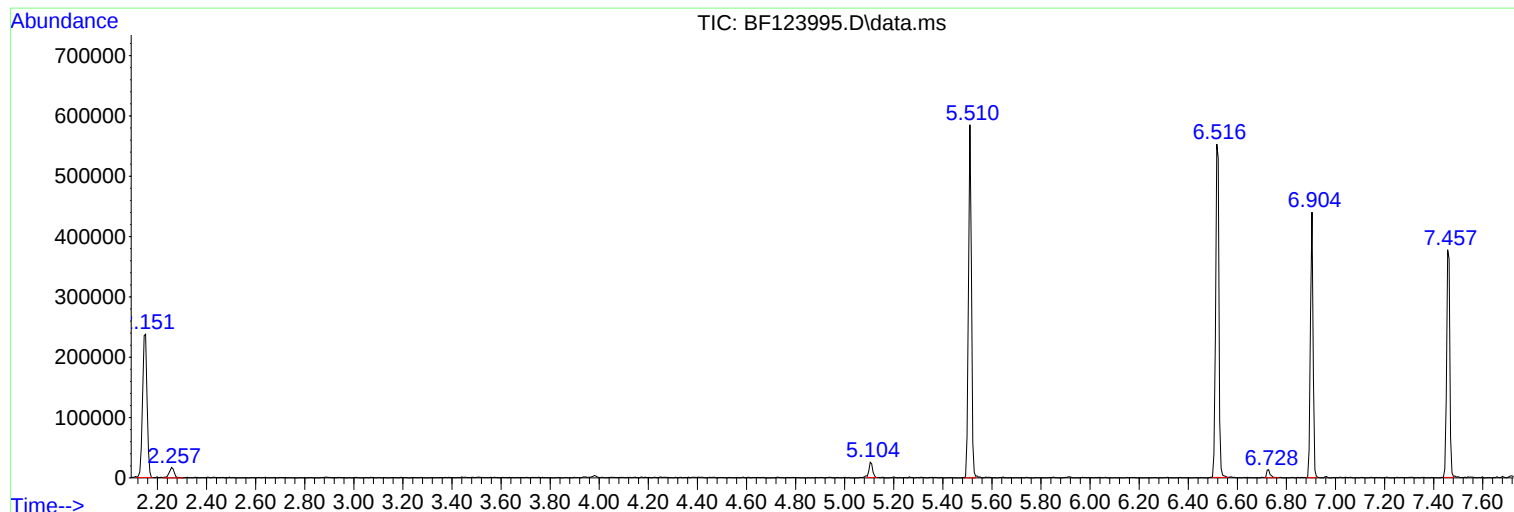
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Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051921\
Data File : BF123995.D
Acq On : 19 May 2021 17:49
Operator : JU/CG
Sample : M2448-01 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
HD-01-051821

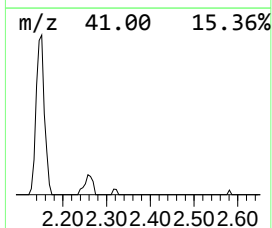
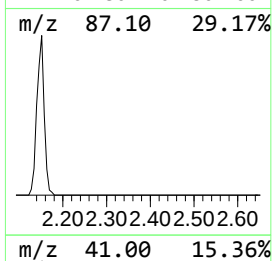
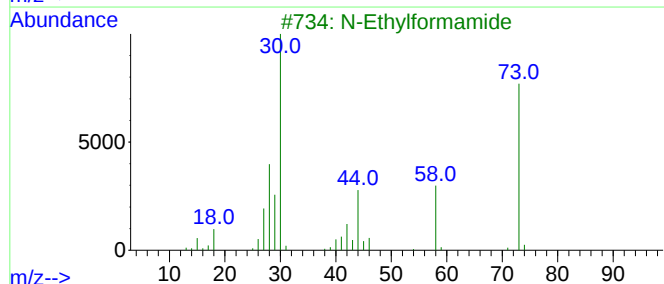
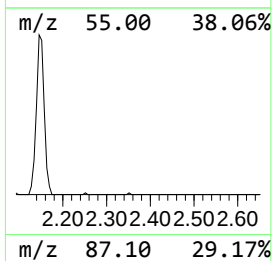
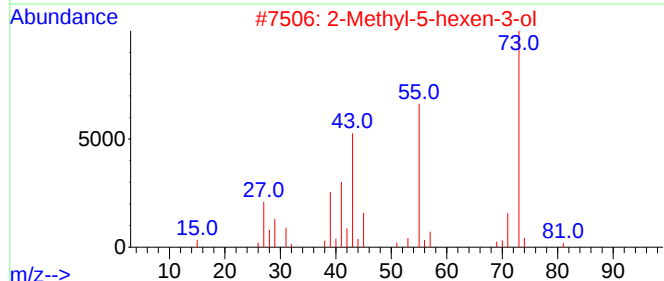
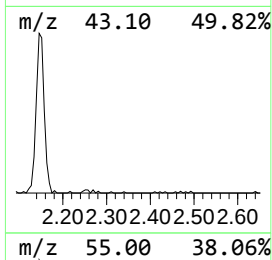
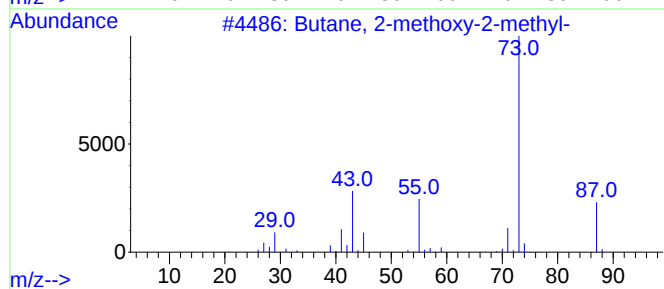
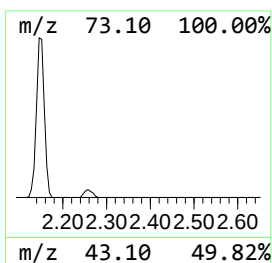
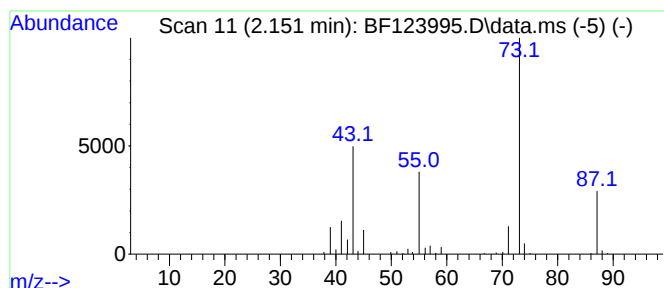
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.151	18.29 ng	303626	1,4-Dichlorobenzene-d4	6.904

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	25
3			N-Ethylformamide	73	C3H7NO	000627-45-2	9
4			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
5			2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9



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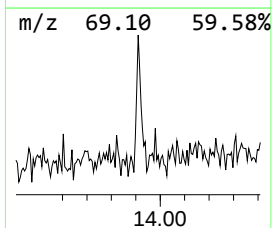
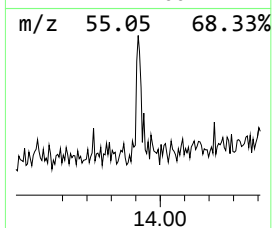
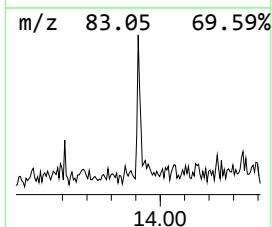
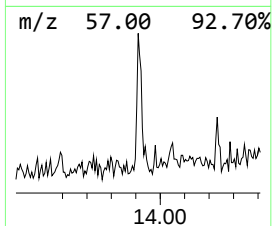
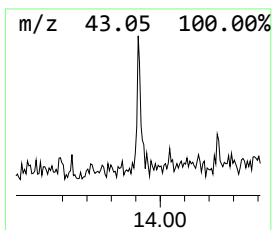
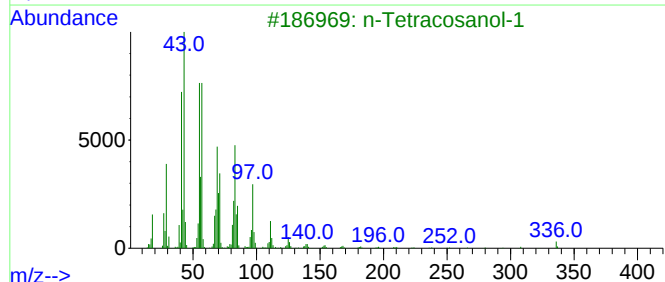
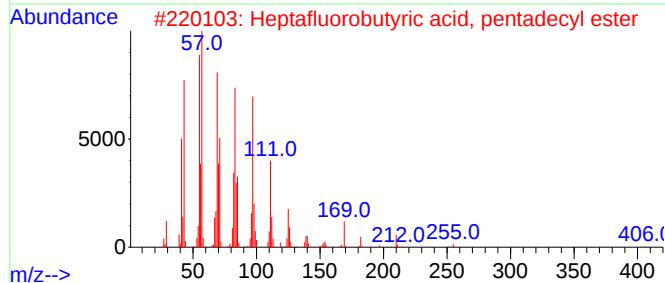
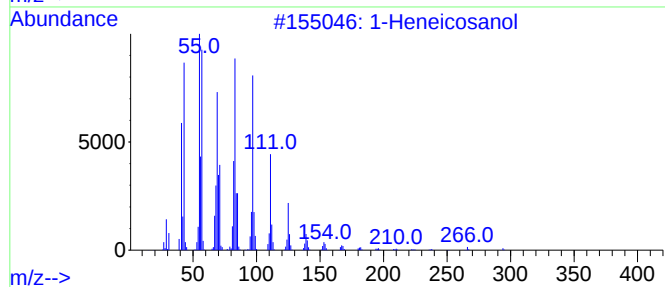
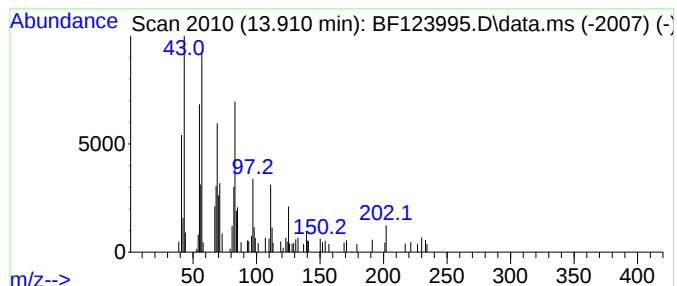
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 1-Heneicosanol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.910	4.36 ng	85628	Chrysene-d12	14.074

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosanol	312	C21H44O	015594-90-8	62
2			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	53
3			n-Tetracosanol-1	354	C24H50O	000506-51-4	53
4			Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	53
5			Heptafluorobutyric acid, n-tetra...	410	C18H29F7O2	007365-36-8	53



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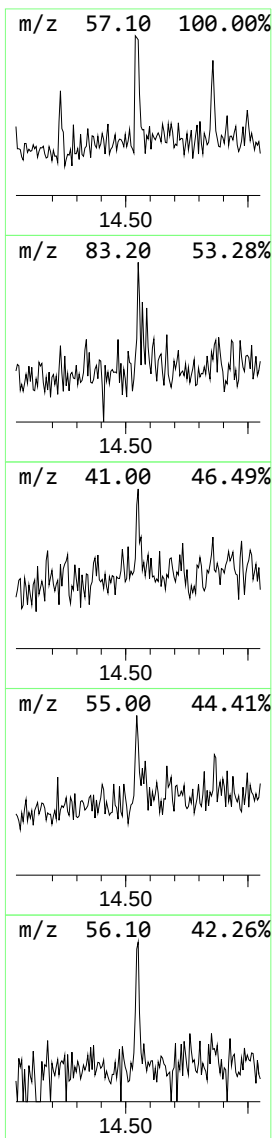
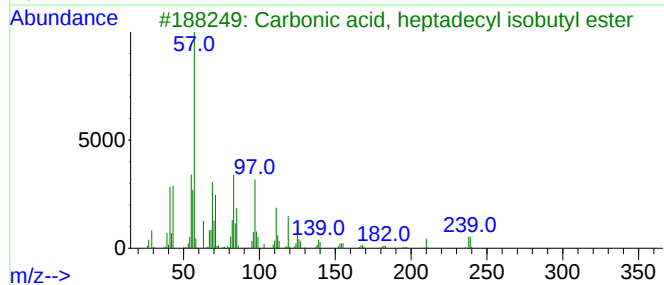
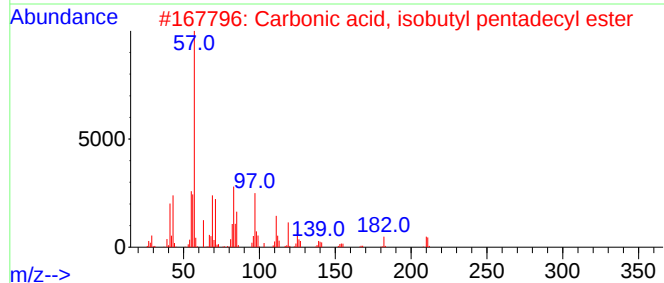
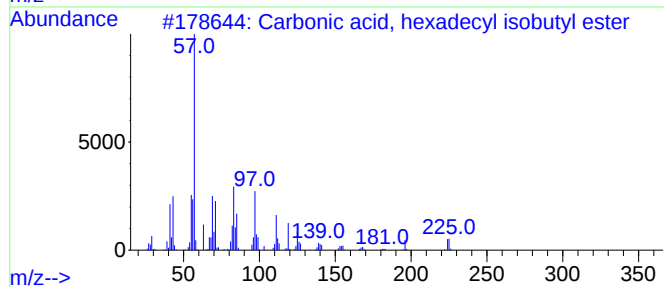
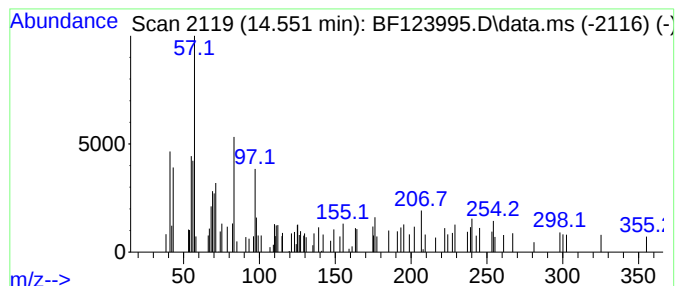
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown14.551 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.551	2.41 ng	47238	Chrysene-d12	14.074

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, hexadecyl isobuty...	342	C21H42O3	1000314-61-3	14
2			Carbonic acid, isobutyl pentadec...	328	C20H40O3	1000314-61-2	14
3			Carbonic acid, heptadecyl isobut...	356	C22H44O3	1000314-61-4	14
4			Hexadecane, 1-(ethenylxy)-	268	C18H36O	000822-28-6	14
5			2,2-Dicyclohexylbutane	222	C16H30	054890-02-7	10



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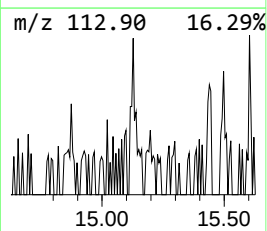
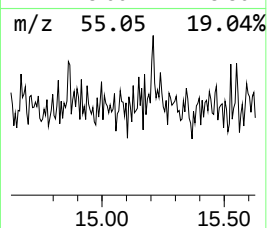
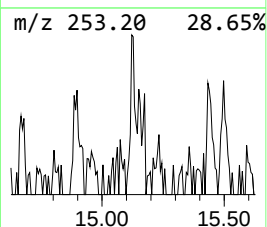
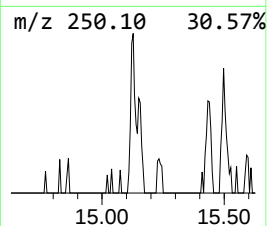
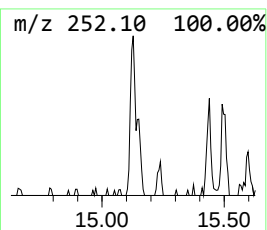
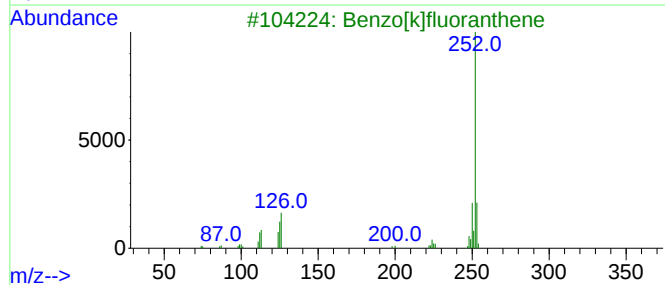
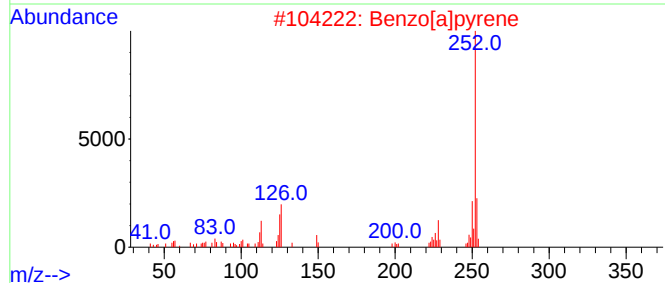
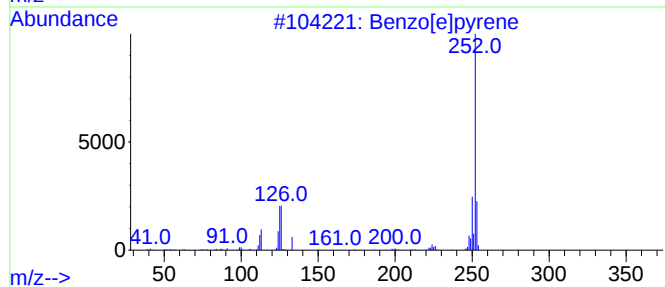
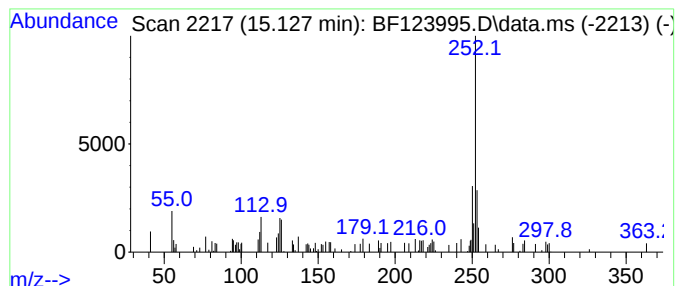
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.127	3.41 ng	49517	Perylene-d12	15.568

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	70
2			Benzo[a]pyrene	252	C20H12	000050-32-8	70
3			Benzo[k]fluoranthene	252	C20H12	000207-08-9	70
4			Perylene	252	C20H12	000198-55-0	58
5			2,1,3-Benzoselenadiazole, 4,6-di...	252	C6H2C12N2Se	090030-99-2	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051921\
Data File : BF123995.D
Acq On : 19 May 2021 17:49
Operator : JU/CG
Sample : M2448-01 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HD-01-051821

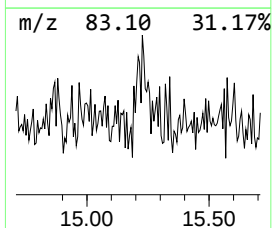
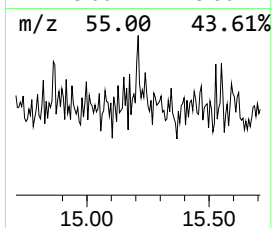
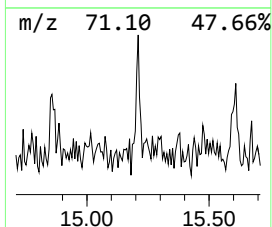
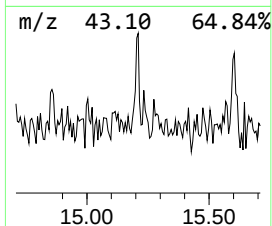
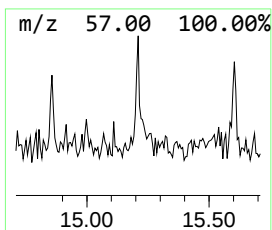
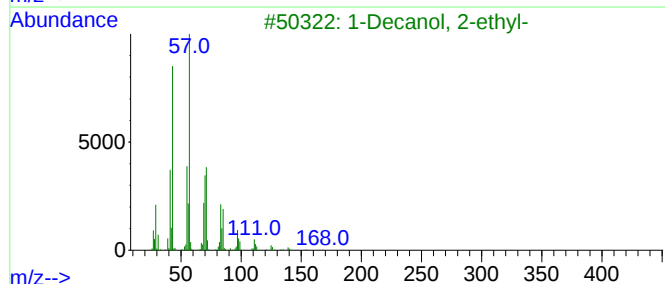
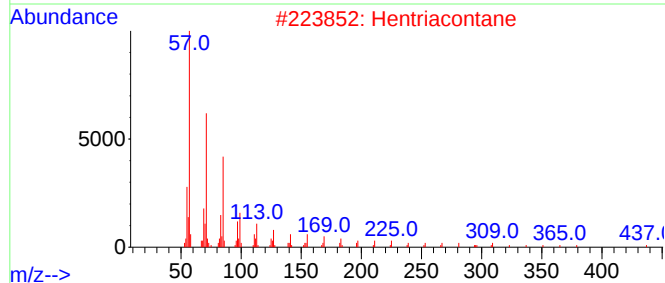
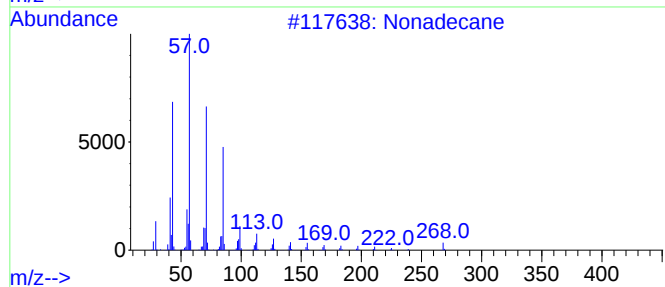
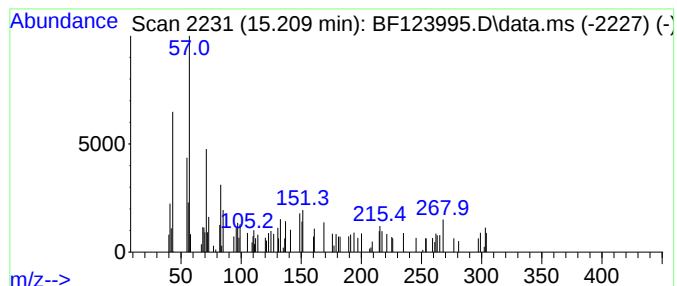
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 unknown15.210 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.210	2.45 ng	35608	Perylene-d12	15.568

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	38
2			Hentriacontane	437	C31H64	000630-04-6	35
3			1-Decanol, 2-ethyl-	186	C12H26O	021078-65-9	27
4			Undecane, 3,5-dimethyl-	184	C13H28	017312-81-1	27
5			Octane, 2,7-dimethyl-	142	C10H22	001072-16-8	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051921\
Data File : BF123995.D
Acq On : 19 May 2021 17:49
Operator : JU/CG
Sample : M2448-01 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HD-01-051821

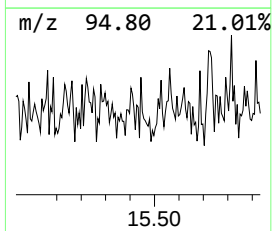
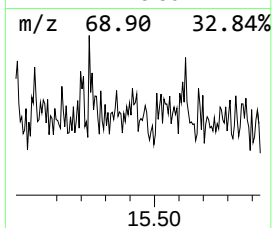
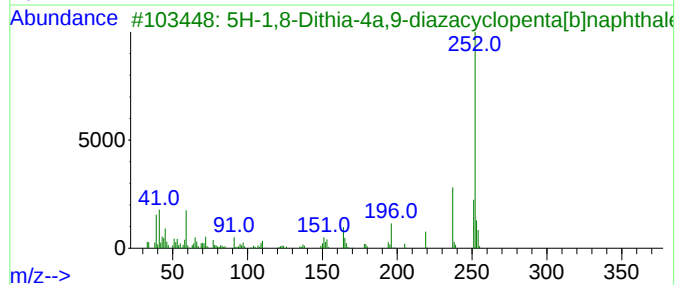
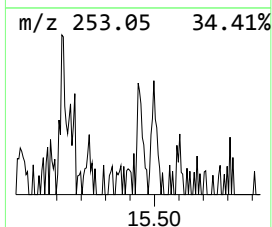
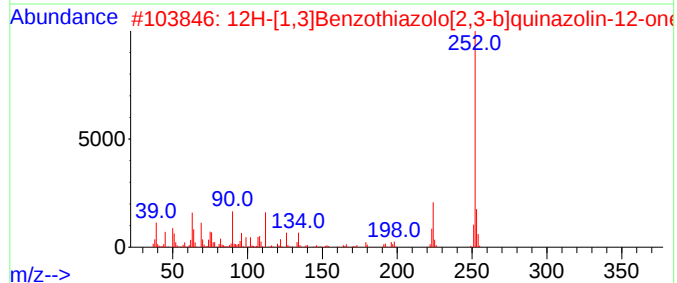
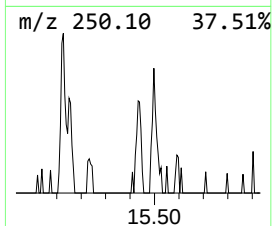
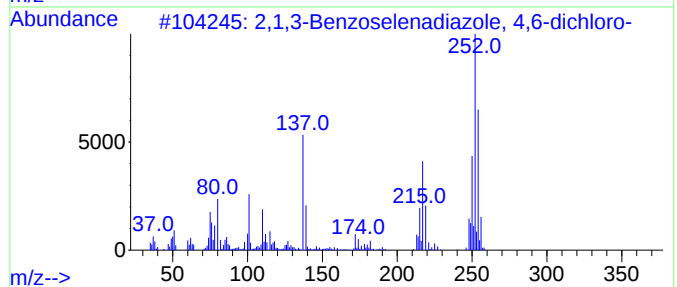
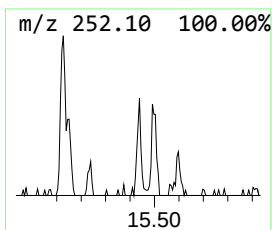
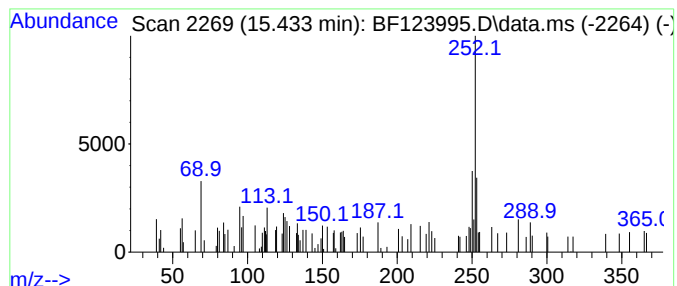
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown15.433 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.433	2.78 ng	40422	Perylene-d12	15.568

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,1,3-Benzoselenadiazole, 4,6-di...	252	C6H2Cl2N2Se	090030-99-2	49		
2	12H-[1,3]Benzothiazolo[2,3-b]qui...	252	C14H8N2OS	1000338-32-6	43		
3	5H-1,8-Dithia-4a,9-diazacyclop...	252	C11H12N2OS2	1000315-96-2	38		
4	Benzo[e]pyrene	252	C20H12	000192-97-2	38		
5	1-Amino-4-(phenylthio)isoquinoline	252	C15H12N2S	019653-19-1	38		



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051921\
Data File : BF123995.D
Acq On : 19 May 2021 17:49
Operator : JU/CG
Sample : M2448-01 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

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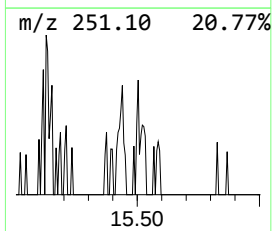
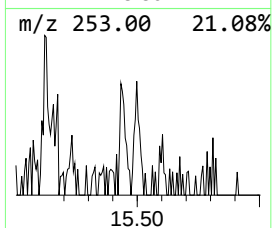
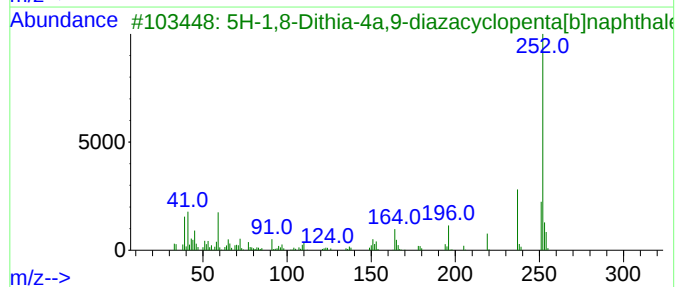
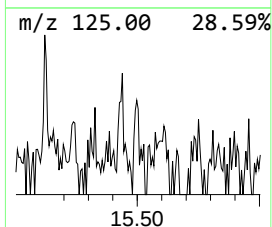
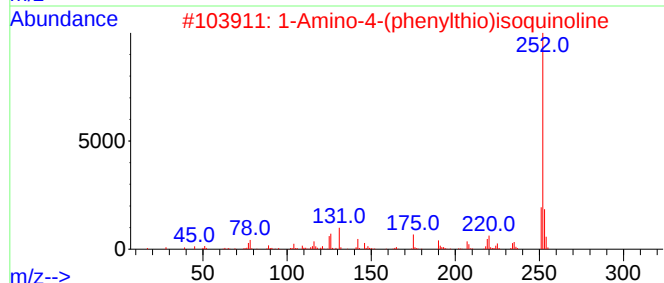
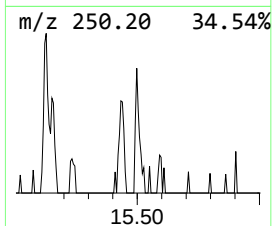
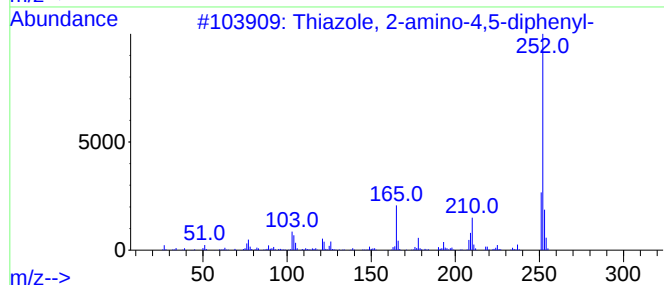
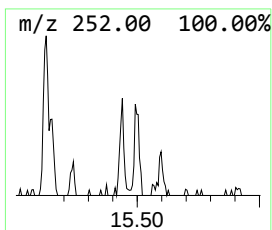
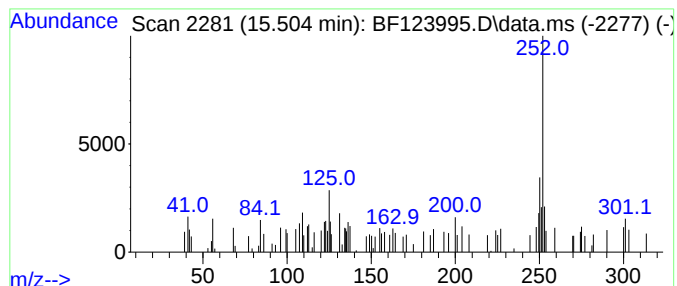
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown15.504 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.504	2.40 ng	34918	Perylene-d12	15.568

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Thiazole, 2-amino-4,5-diphenyl-	252	C15H12N2S	006318-74-7	46
2			1-Amino-4-(phenylthio)isoquinoline	252	C15H12N2S	019653-19-1	46
3			5H-1,8-Dithia-4a,9-diazacyclopenta[b]naphthalen-2-amine	252	C11H12N2OS2	1000315-96-2	43
4			1H-1,3-Benzimidazole, 2-tricyclopropyl-	252	C17H20N2	1000319-80-9	43
5			N-(3-Nitro-thiobenzoyl)-morpholine	252	C11H12N2O3S	050903-05-4	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051921\
Data File : BF123995.D
Acq On : 19 May 2021 17:49
Operator : JU/CG
Sample : M2448-01 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

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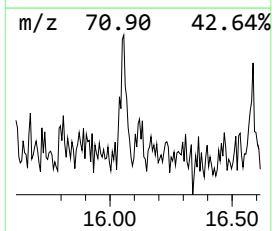
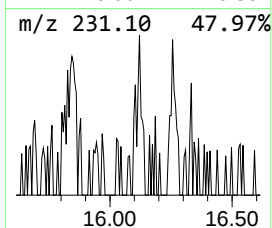
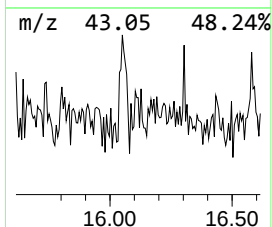
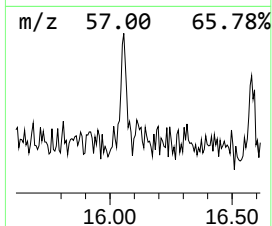
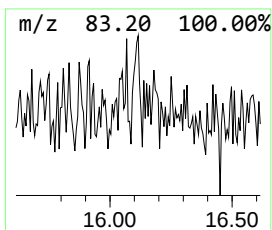
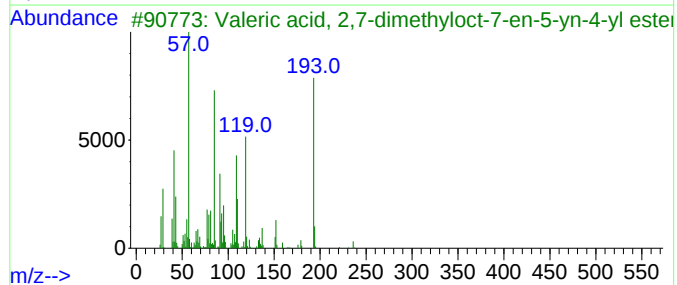
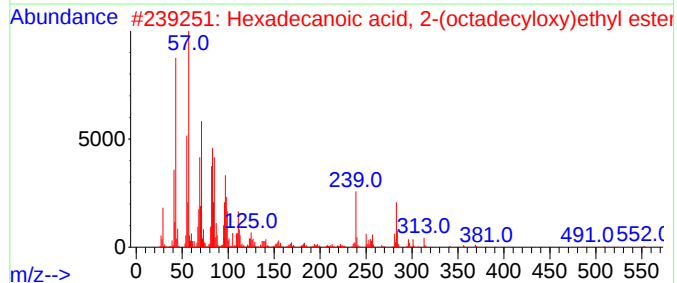
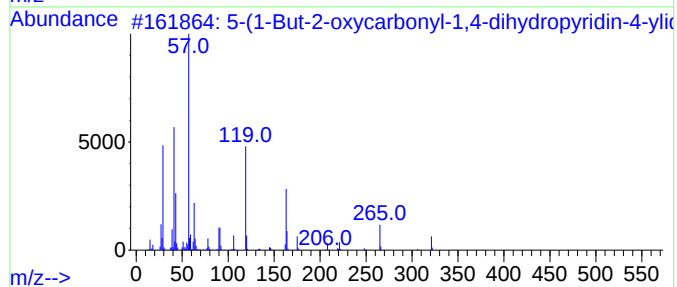
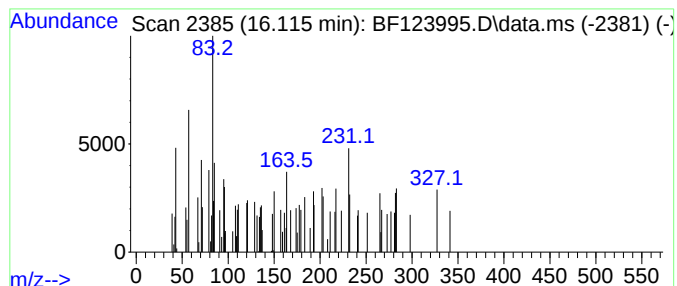
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown16.115 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.115	2.70 ng	39154	Perylene-d12	15.568

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-(1-But-2-oxycarbonyl-1,4-dihyd...	321	C16H19N06	142131-83-7	9		
2	Hexadecanoic acid, 2-(octadecylo...	553	C36H72O3	029899-13-6	7		
3	Valeric acid, 2,7-dimethyloct-7-...	236	C15H24O2	1000292-48-9	7		
4	Acetyl 2,4-di-O-acetyl-3,6-dideo...	274	C12H18O7	1000101-82-5	2		
5	4-tert-Butylsulfanyl-6-fluoro-2-...	287	C15H14FN3S	1000275-85-6	1		



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051921\
Data File : BF123995.D
Acq On : 19 May 2021 17:49
Operator : JU/CG
Sample : M2448-01 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
HD-01-051821

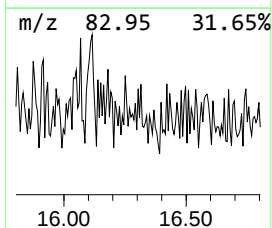
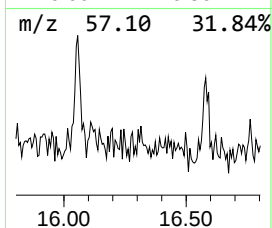
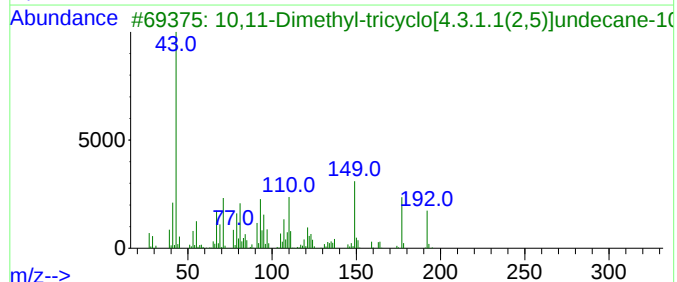
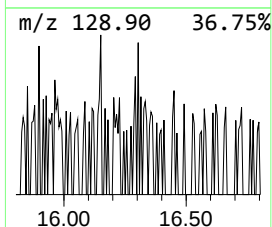
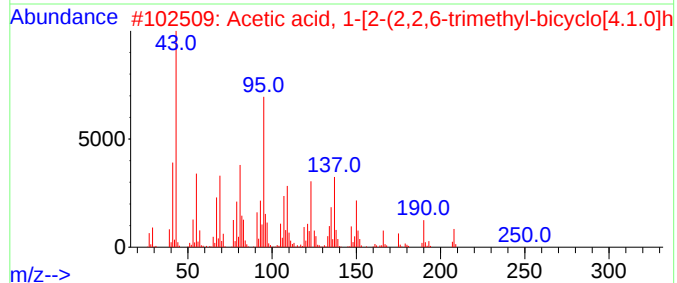
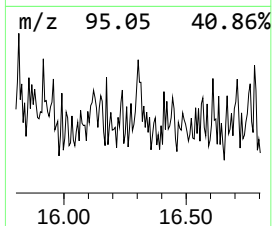
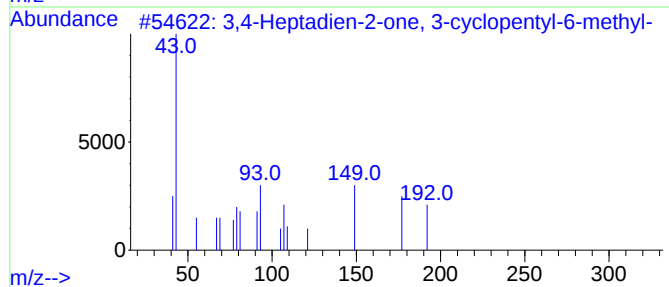
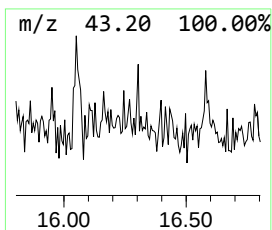
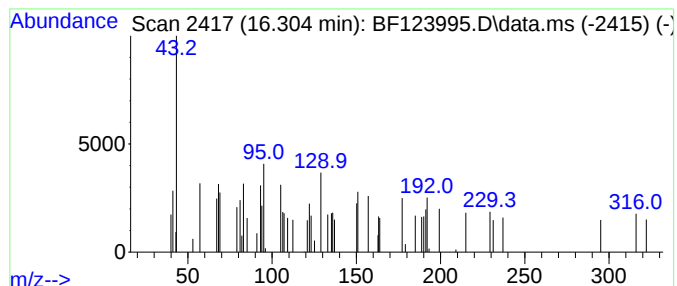
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown16.304 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.304	2.42 ng	35119	Perylene-d12	15.568

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,4-Heptadien-2-one, 3-cyclopent...	192	C13H20O	063922-50-9	15
2			Acetic acid, 1-[2-(2,2,6-trimeth...	250	C16H26O2	077143-23-8	10
3			10,11-Dimethyl-tricyclo[4.3.1.1(...	210	C13H22O2	174226-41-6	10
4			3-(3,4-Dimethoxycinnamoyl)-4-hyd...	316	C17H16O6	1000225-16-8	9
5			Benzaldehyde, 4-methoxy-3-(2-pro...	192	C11H12O3	018075-40-6	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051921\
Data File : BF123995.D
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Misc :
ALS Vial : 13 Sample Multiplier: 1

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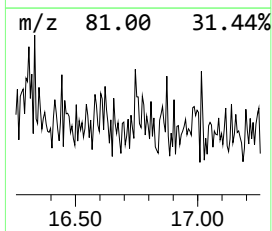
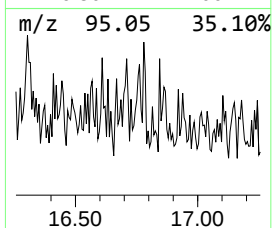
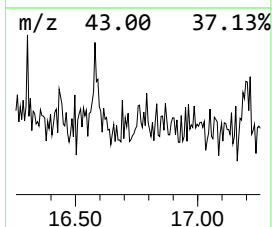
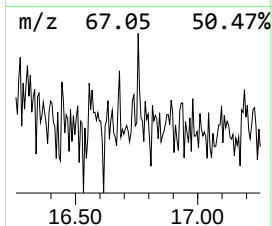
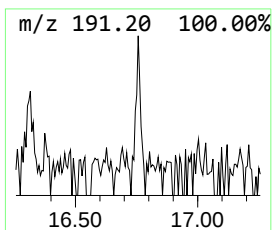
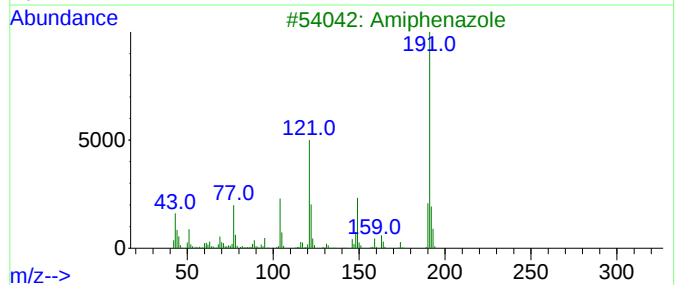
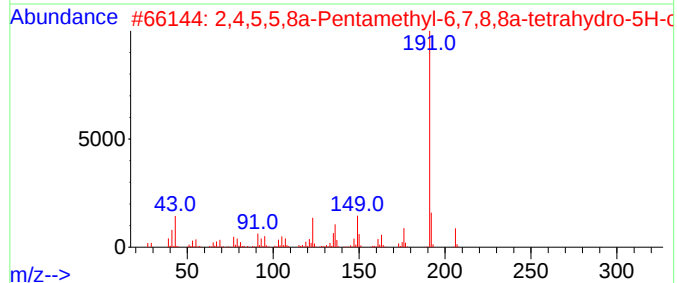
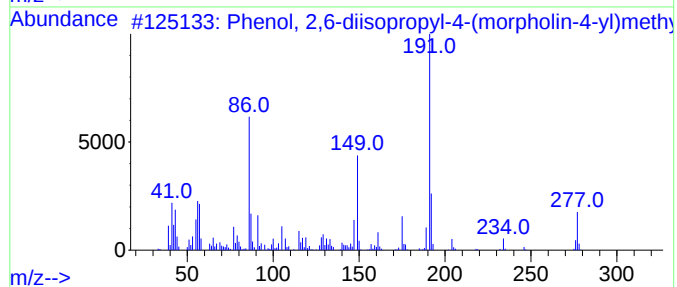
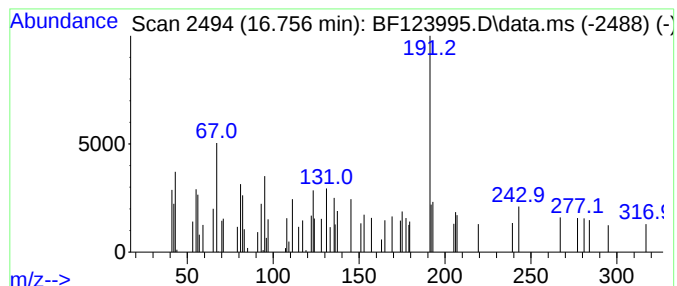
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown16.756 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.756	3.04 ng	44067	Perylene-d12	15.568

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,6-diisopropyl-4-(morph...	277	C17H27NO2	1000304-55-0	46
2			2,4,5,5,8a-Pentamethyl-6,7,8,8a-...	206	C14H22O	1000195-40-9	43
3			Amiphenazole	191	C9H9N3S	000490-55-1	38
4			Silane, 1,3-decadiynyltrimethyl-	206	C13H22Si	084751-17-7	38
5			2-Anthracenyloxirane	220	C16H12O	052643-86-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051921\
Data File : BF123995.D
Acq On : 19 May 2021 17:49
Operator : JU/CG
Sample : M2448-01 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HD-01-051821

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	2.151	18.3	ng	303626	1	6.904	331985	20.0
1-Heneicosanol	13.910	4.4	ng	85628	5	14.074	392371	20.0
unknown14.551	14.551	2.4	ng	47238	5	14.074	392371	20.0
Benzo[e]pyrene	15.127	3.4	ng	49517	6	15.568	290388	20.0
unknown15.210	15.210	2.5	ng	35608	6	15.568	290388	20.0
unknown15.433	15.433	2.8	ng	40422	6	15.568	290388	20.0
unknown15.504	15.504	2.4	ng	34918	6	15.568	290388	20.0
unknown16.115	16.115	2.7	ng	39154	6	15.568	290388	20.0
unknown16.304	16.304	2.4	ng	35119	6	15.568	290388	20.0
unknown16.756	16.756	3.0	ng	44067	6	15.568	290388	20.0