

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102320\
 Data File : BF122100.D
 Acq On : 23 Oct 2020 15:46
 Operator : JU/CG
 Sample : L4488-05
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 202

Integration Parameters: rteint.p

Integrator: RTE

Smothing : 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-BF101220.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF122100.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.293	372	375	391	rBV2	20981	47951	1.79%	0.209%
2	4.940	481	485	495	rBV	28056	49352	1.84%	0.215%
3	5.345	551	554	580	rBV	1692842	1902198	70.84%	8.271%
4	6.375	725	729	757	rBV	2219005	2277338	84.81%	9.903%
5	6.581	762	764	783	rVV	20773	61621	2.29%	0.268%
6	6.722	785	788	799	rVB	707221	588384	21.91%	2.558%
7	7.292	881	885	907	rBV	1603470	1628241	60.64%	7.080%
8	8.010	1003	1007	1016	rBV	763405	794041	29.57%	3.453%
9	9.080	1184	1189	1199	rBV	2855108	2685297	100.00%	11.676%
10	9.480	1254	1257	1264	rVB	387666	311639	11.61%	1.355%
11	9.757	1300	1304	1308	rBV	978205	914843	34.07%	3.978%
12	9.792	1308	1310	1315	rVB	39767	36035	1.34%	0.157%
13	9.969	1336	1340	1344	rVB2	27032	29452	1.10%	0.128%
14	10.163	1369	1373	1375	rBV2	29801	34479	1.28%	0.150%
15	10.310	1392	1398	1403	rBV3	30173	48778	1.82%	0.212%
16	10.404	1410	1414	1419	rVB3	21056	33021	1.23%	0.144%
17	10.551	1436	1439	1449	rBV	1012588	951419	35.43%	4.137%
18	10.669	1455	1459	1462	rBV	70613	71692	2.67%	0.312%
19	10.986	1507	1513	1515	rBV3	28338	40699	1.52%	0.177%
20	11.104	1529	1533	1535	rBV2	26780	36406	1.36%	0.158%
21	11.139	1535	1539	1546	rVB3	68864	100854	3.76%	0.439%
22	11.245	1553	1557	1560	rBV	1052869	984634	36.67%	4.281%
23	11.269	1560	1561	1565	rVV	424959	306537	11.42%	1.333%
24	11.327	1567	1571	1574	rVB	94505	90215	3.36%	0.392%
25	11.492	1594	1599	1605	rBV5	47228	103420	3.85%	0.450%
26	11.751	1640	1643	1645	rBV	71215	79158	2.95%	0.344%
27	11.780	1645	1648	1654	rVB	244202	238708	8.89%	1.038%
28	11.827	1654	1656	1659	rVB	46551	32789	1.22%	0.143%
29	11.863	1659	1662	1664	rBV	84798	87053	3.24%	0.379%
30	12.045	1690	1693	1697	rVB	52323	45090	1.68%	0.196%
31	12.092	1698	1701	1703	rVV3	38498	32438	1.21%	0.141%
32	12.192	1713	1718	1720	rBV4	34001	51897	1.93%	0.226%
33	12.298	1733	1736	1738	rBV	58122	54527	2.03%	0.237%
34	12.321	1738	1740	1741	rVV2	52025	47376	1.76%	0.206%
35	12.339	1741	1743	1745	rVV2	72445	62420	2.32%	0.271%
36	12.398	1748	1753	1755	rBV2	37549	47286	1.76%	0.206%

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Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
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37	12.463	1760	1764	1767	rBV	617571	524184	19.52%	2.279%
38	12.563	1773	1781	1787	rBV	288828	425303	15.84%	1.849%
39	12.616	1787	1790	1793	rVV	65708	93366	3.48%	0.406%
40	12.692	1797	1803	1809	rVV2	470384	583565	21.73%	2.538%
41	12.745	1809	1812	1816	rVB2	44313	47778	1.78%	0.208%
42	12.833	1821	1827	1830	rVV	2723213	2530026	94.22%	11.001%
43	12.863	1830	1832	1836	rVB2	26692	31627	1.18%	0.138%
44	12.910	1836	1840	1844	rBV2	42261	73945	2.75%	0.322%
45	12.963	1847	1849	1852	rVB	59320	45192	1.68%	0.197%
46	12.998	1852	1855	1856	rBV2	25195	29924	1.11%	0.130%
47	13.021	1856	1859	1862	rVB	60433	60550	2.25%	0.263%
48	13.057	1862	1865	1868	rBV3	20643	30256	1.13%	0.132%
49	13.121	1872	1876	1880	rBV2	54819	68108	2.54%	0.296%
50	13.251	1895	1898	1901	rBV2	29831	33856	1.26%	0.147%
51	13.363	1915	1917	1920	rVB	43588	32648	1.22%	0.142%
52	13.539	1944	1947	1950	rBV	51965	48507	1.81%	0.211%
53	13.645	1962	1965	1967	rBV2	36805	40618	1.51%	0.177%
54	13.663	1967	1968	1971	rVB	46268	32869	1.22%	0.143%
55	13.692	1971	1973	1975	rBV	31034	30307	1.13%	0.132%
56	13.751	1979	1983	1989	rVB	168138	262112	9.76%	1.140%
57	13.862	1998	2002	2003	rBV	68607	78883	2.94%	0.343%
58	13.892	2003	2007	2009	rVV2	786695	910658	33.91%	3.960%
59	13.915	2009	2011	2015	rVB	208080	187927	7.00%	0.817%
60	14.039	2029	2032	2034	rBV2	26313	28959	1.08%	0.126%
61	14.280	2071	2073	2076	rVB	46762	40042	1.49%	0.174%
62	14.345	2081	2084	2088	rVV2	35167	43700	1.63%	0.190%
63	14.504	2108	2111	2116	rBV	58525	66062	2.46%	0.287%
64	14.709	2143	2146	2152	rVB	183453	188275	7.01%	0.819%
65	14.774	2154	2157	2161	rVB2	28698	30490	1.14%	0.133%
66	14.921	2178	2182	2184	rBV	162211	220567	8.21%	0.959%
67	15.021	2197	2199	2206	rVB	39087	49988	1.86%	0.217%
68	15.209	2228	2231	2236	rBV	89979	94371	3.51%	0.410%
69	15.268	2238	2241	2245	rVB	139127	146333	5.45%	0.636%
70	15.327	2246	2251	2255	rBV	535690	653571	24.34%	2.842%
71	15.721	2315	2318	2322	rBV	39755	58343	2.17%	0.254%
72	16.745	2486	2492	2499	rBV	81980	165877	6.18%	0.721%
73	17.168	2560	2564	2576	rVB	43099	101391	3.78%	0.441%

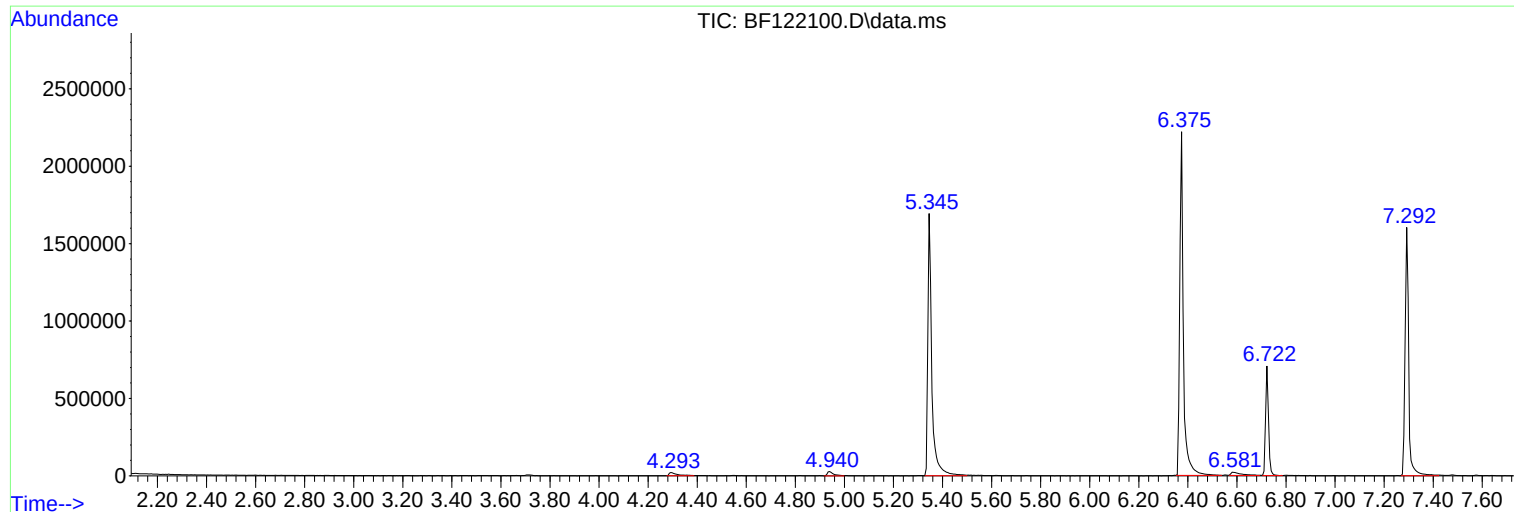
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Instrument :
BNA_F
ClientSampled :
202

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.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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Instrument :
BNA_F
ClientSampleId :
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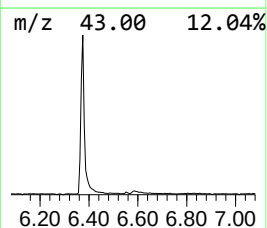
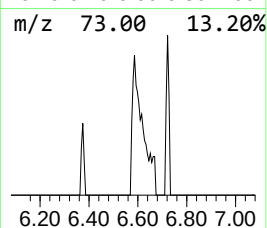
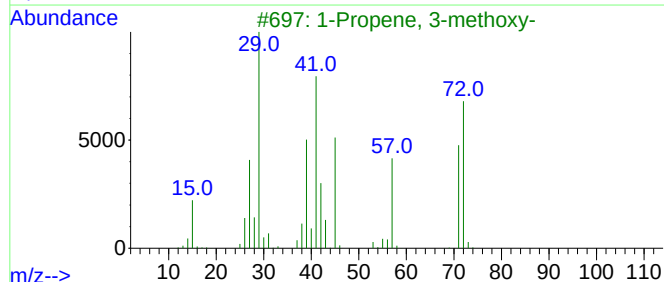
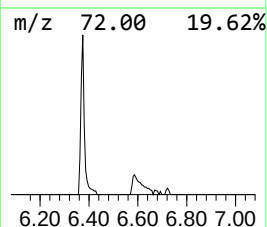
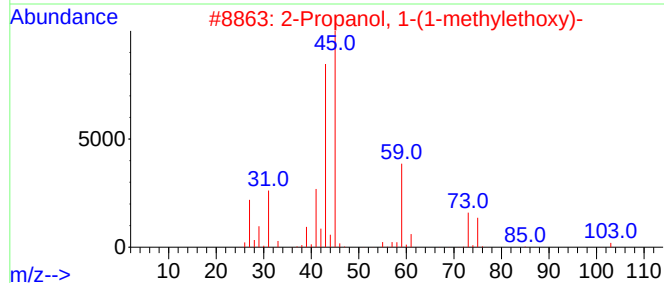
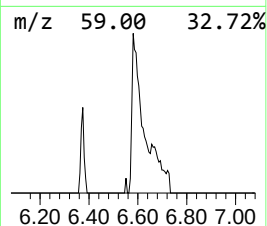
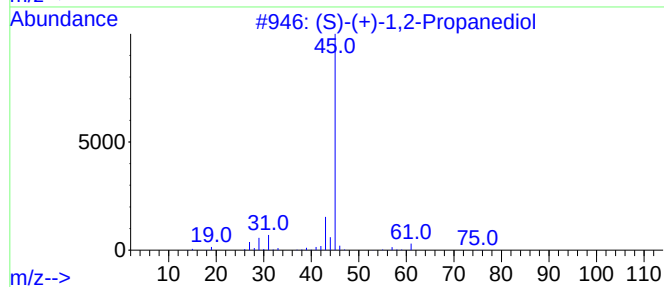
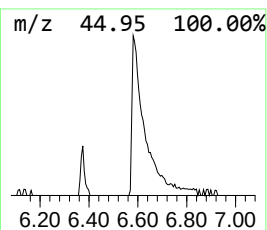
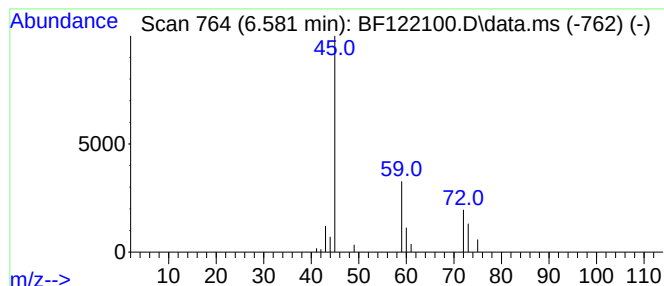
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.58 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.581	2.09 ng	61621	1,4-Dichlorobenzene-d4	6.722

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(S)-(+)-1,2-Propanediol			76	C3H8O2	004254-15-3	9
2	2-Propanol, 1-(1-methylethoxy)-			118	C6H14O2	003944-36-3	9
3	1-Propene, 3-methoxy-			72	C4H8O	000627-40-7	4
4	Methane, nitroso-			45	CH3NO	000865-40-7	4
5	Hydrazine, 2-propenyl-			72	C3H8N2	007422-78-8	4



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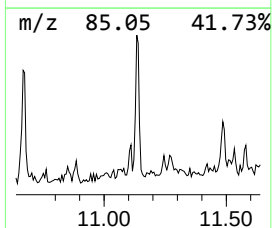
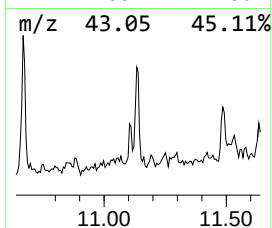
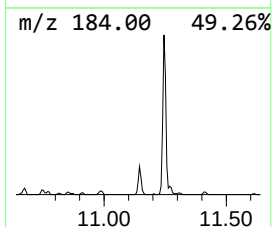
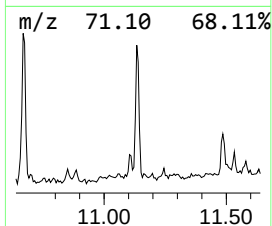
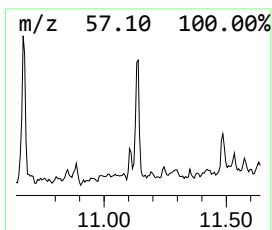
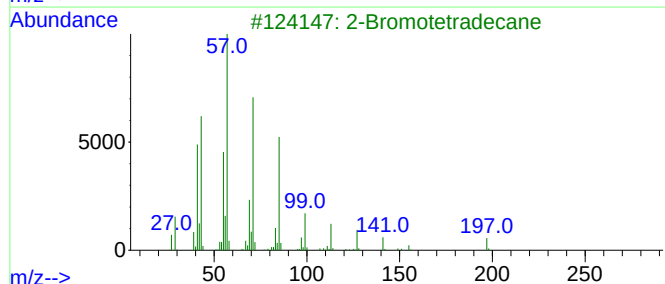
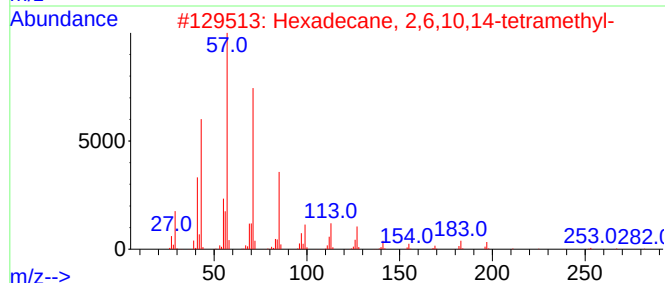
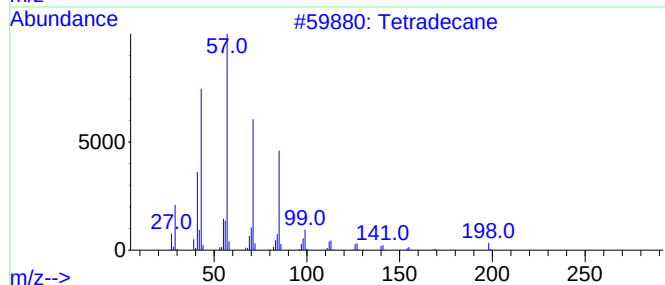
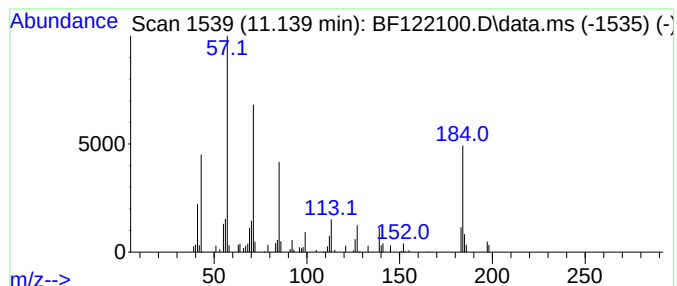
Instrument :
BNA_F
ClientSampleId :
202

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Tetradecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.139	2.05 ng	100854	Phenanthrene-d10		11.245	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tetradecane		198	C14H30	000629-59-4	53
2	Hexadecane, 2,6,10,14-tetramethyl-		282	C20H42	000638-36-8	52
3	2-Bromotetradecane		276	C14H29Br	074036-95-6	47
4	Sulfurous acid, 2-ethylhexyl hex...		278	C14H30O3S	1000309-20-2	46
5	3-Ethyl-3-methylheptane		142	C10H22	017302-01-1	43



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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
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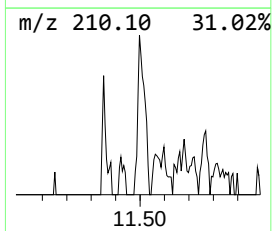
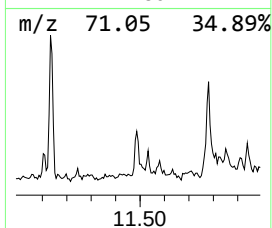
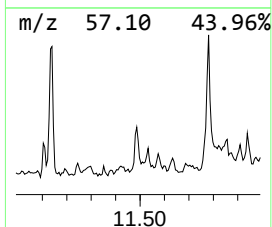
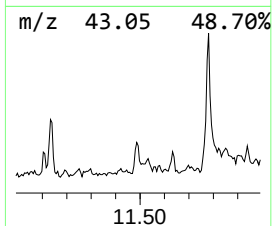
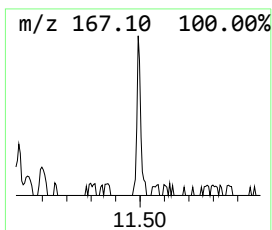
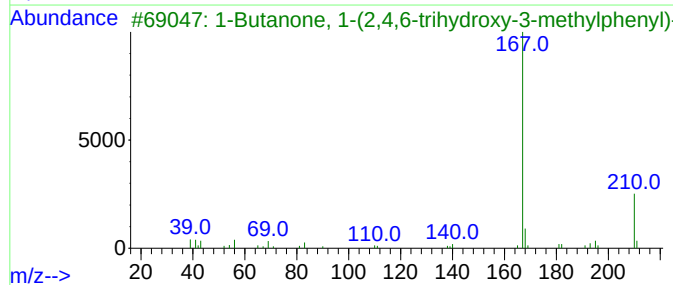
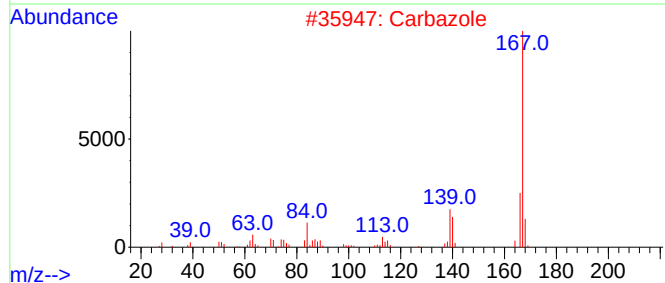
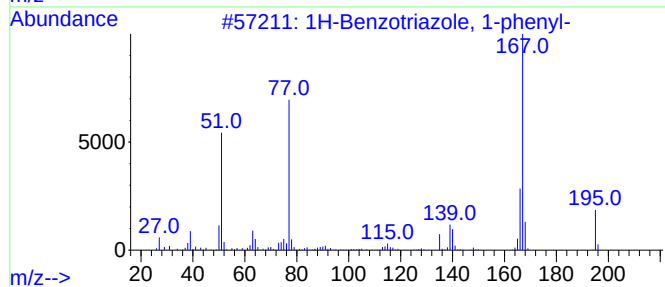
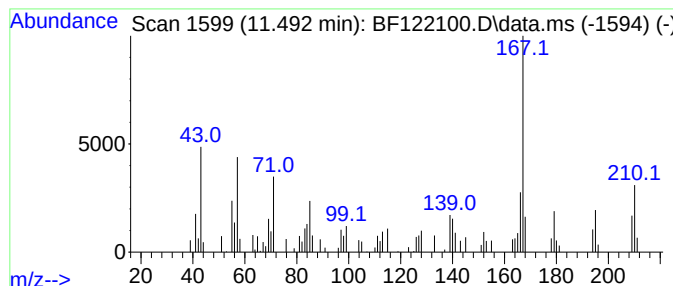
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 1H-Benzotriazole, 1-phenyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.492	2.10 ng	103420	Phenanthrene-d10	11.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Benzotriazole, 1-phenyl-	195	C12H9N3	000883-39-6	52
2			Carbazole	167	C12H9N	000086-74-8	45
3			1-Butanone, 1-(2,4,6-trihydroxy-...	210	C11H14O4	001509-06-4	43
4			2-Propanone, 1,1-diphenyl-	210	C15H14O	000781-35-1	43
5			5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	43



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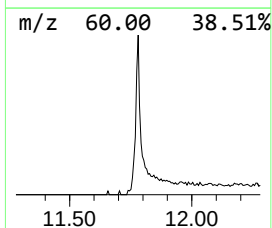
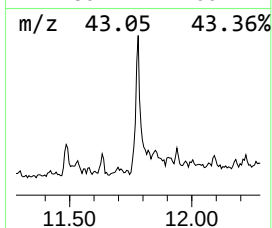
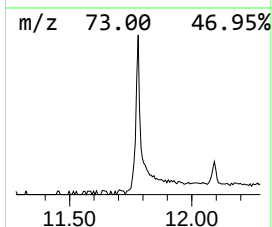
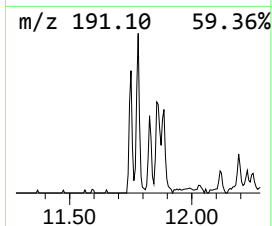
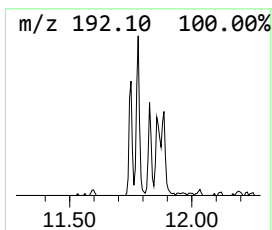
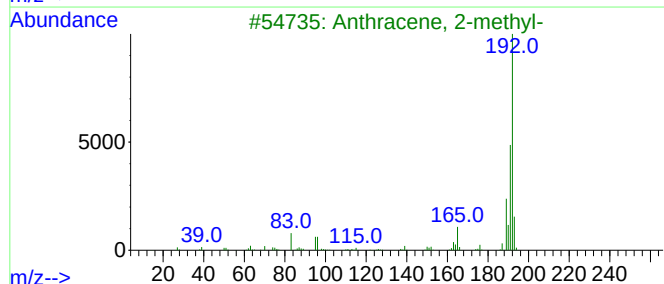
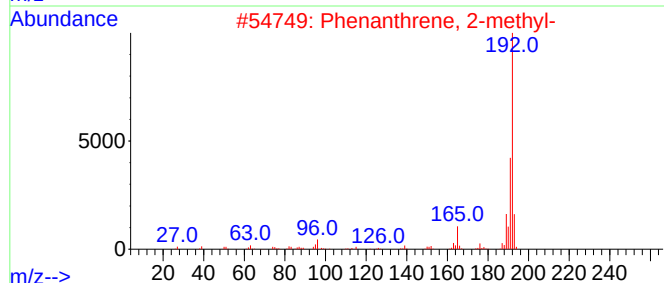
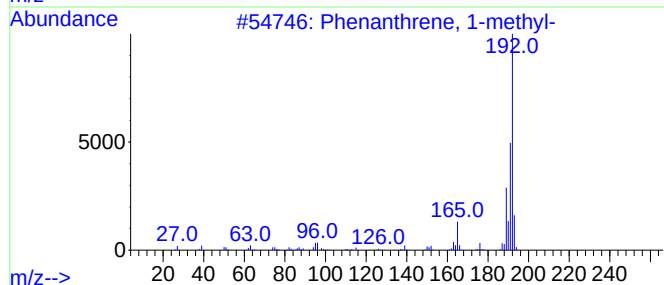
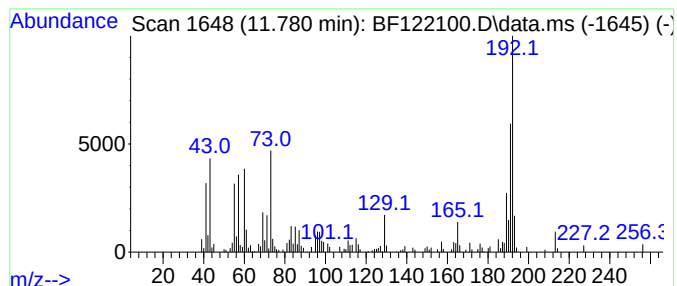
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Phenanthrene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.780	4.85 ng	238708	Phenanthrene-d10	11.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
4			Anthracene, 1-methyl-	192	C15H12	000610-48-0	95
5			Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102320\
Data File : BF122100.D
Acq On : 23 Oct 2020 15:46
Operator : JU/CG
Sample : L4488-05
Misc :
ALS Vial : 6 Sample Multiplier: 1

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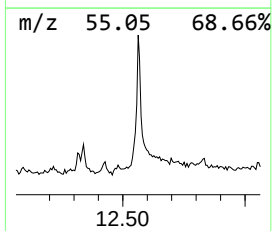
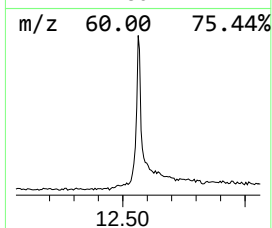
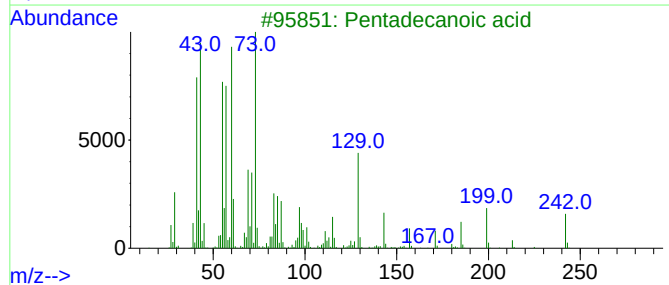
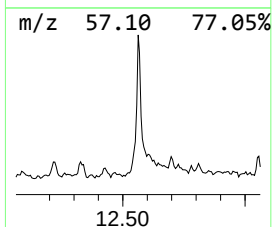
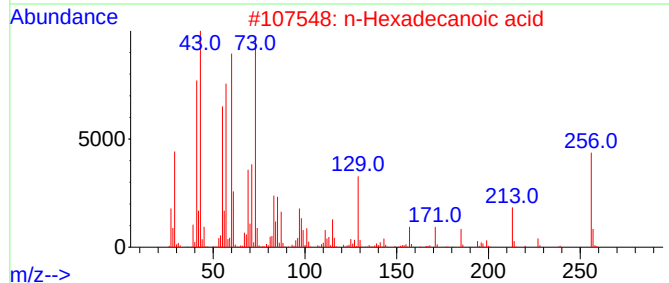
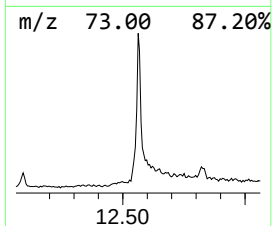
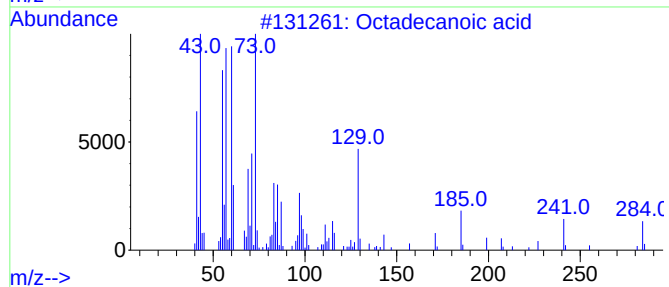
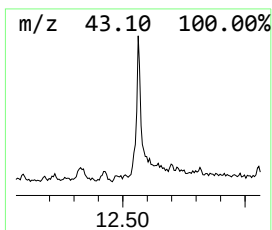
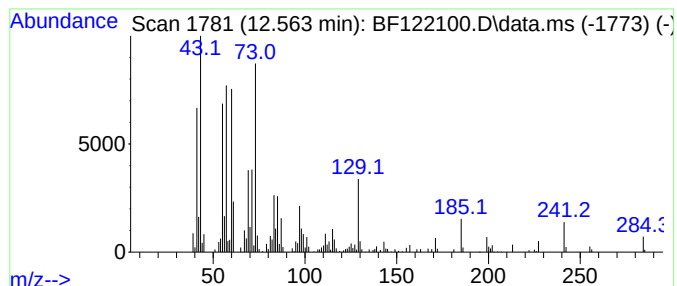
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Octadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.563	8.64 ng	425303	Phenanthrene-d10	11.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	93
4			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	91
5			n-Decanoic acid	172	C10H20O2	000334-48-5	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102320\
Data File : BF122100.D
Acq On : 23 Oct 2020 15:46
Operator : JU/CG
Sample : L4488-05
Misc :
ALS Vial : 6 Sample Multiplier: 1

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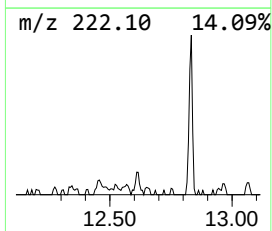
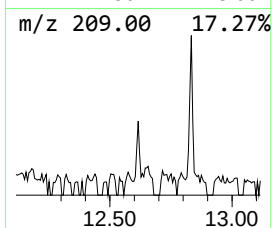
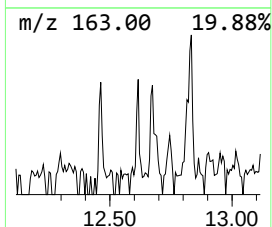
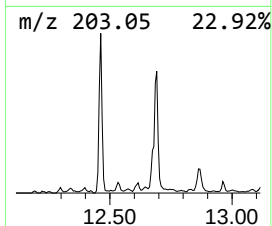
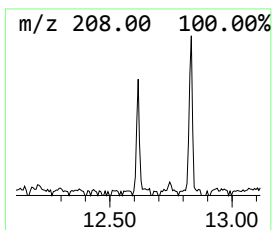
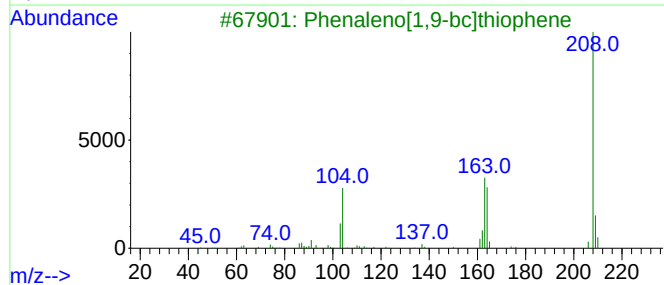
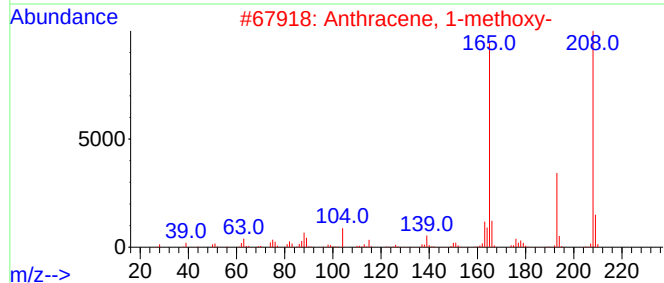
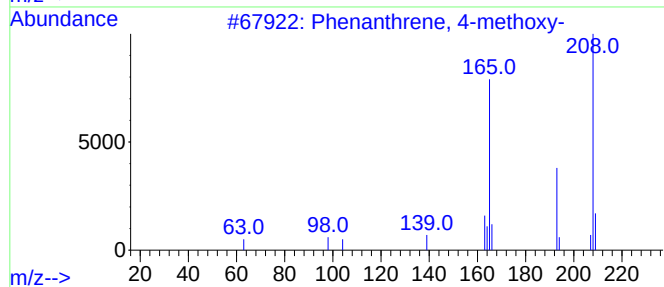
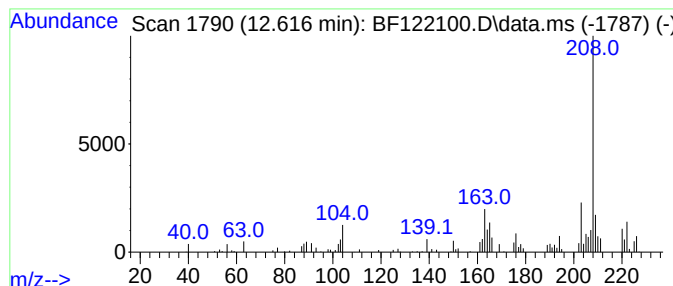
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Phenanthrene, 4-methoxy- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.616	2.05 ng	93366	Chrysene-d12	13.892

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 4-methoxy-	208	C15H12O	015638-06-9	58
2			Anthracene, 1-methoxy-	208	C15H12O	054458-84-3	53
3			Phenalen[1,9-bc]thiophene	208	C14H8S	079965-99-4	50
4			Benzo[c]cinoline, 4,7-dimethyl-	208	C14H12N2	020684-54-2	49
5			6-Methylquinoxaline-2,3-dithiol	208	C9H8N2S2	1000311-61-9	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102320\
Data File : BF122100.D
Acq On : 23 Oct 2020 15:46
Operator : JU/CG
Sample : L4488-05
Misc :
ALS Vial : 6 Sample Multiplier: 1

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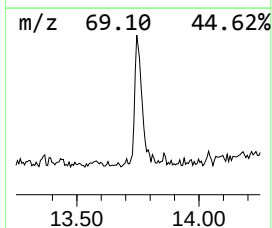
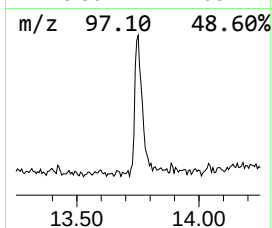
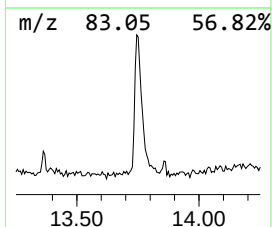
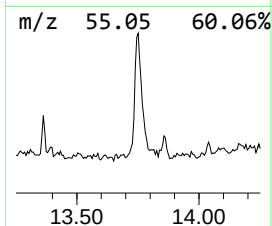
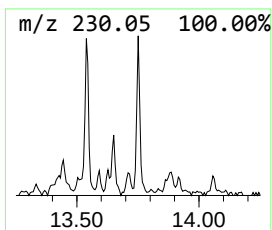
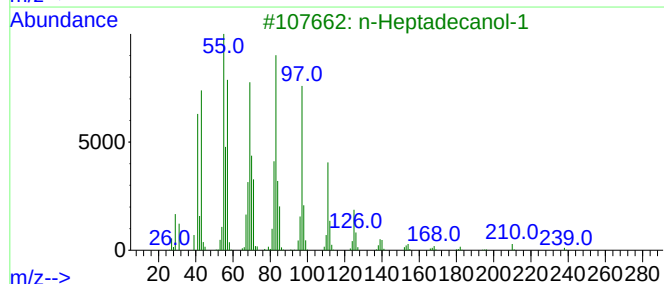
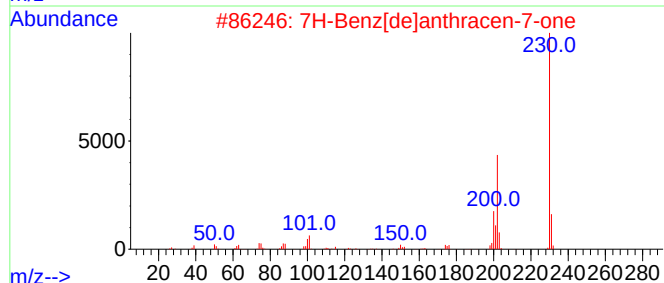
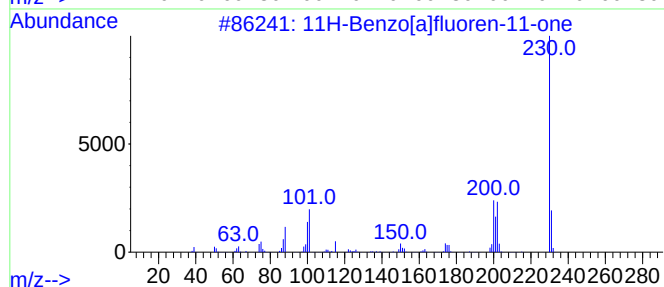
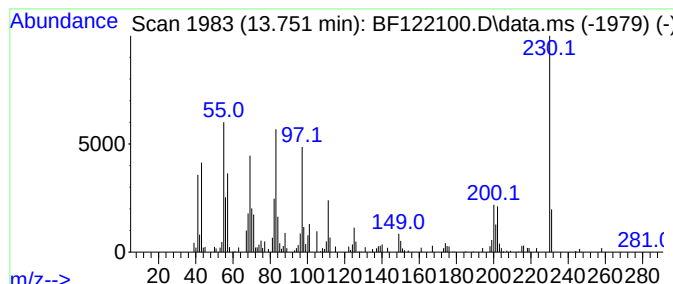
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 11H-Benzo[a]fluoren-11-one Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.751	5.76 ng	262112	Chrysene-d12	13.892

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	95
2			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	50
3			n-Heptadecanol-1	256	C17H36O	001454-85-9	42
4			n-Nonadecanol-1	284	C19H40O	001454-84-8	42
5			6H-Benz[de]anthracen-6-one	230	C17H10O	080252-14-8	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102320\
Data File : BF122100.D
Acq On : 23 Oct 2020 15:46
Operator : JU/CG
Sample : L4488-05
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
202

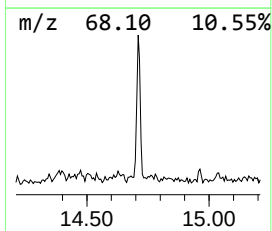
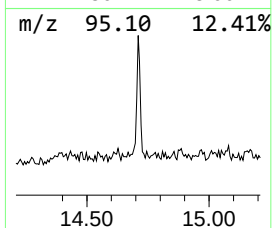
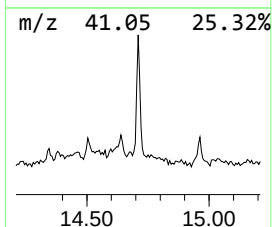
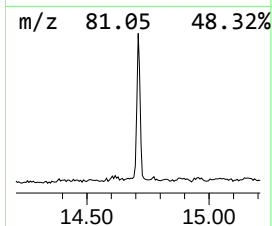
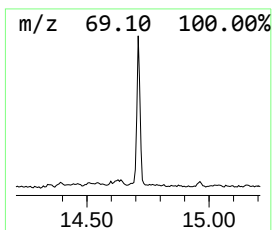
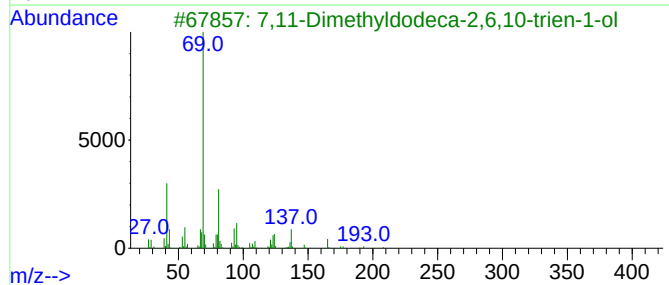
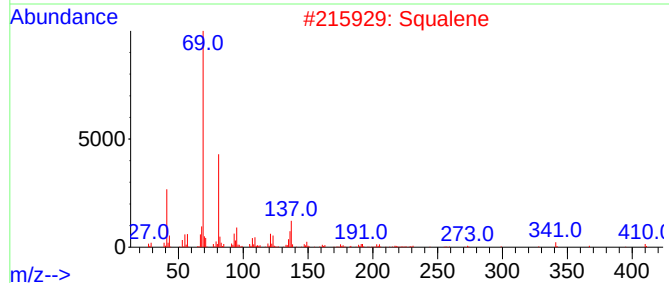
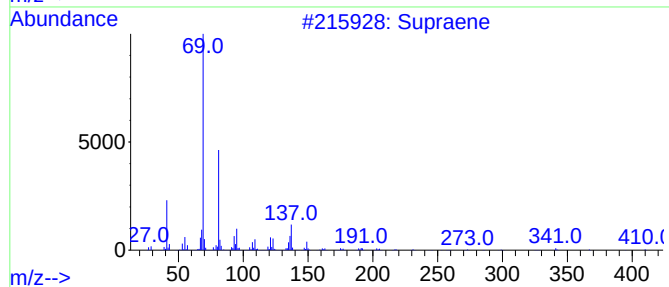
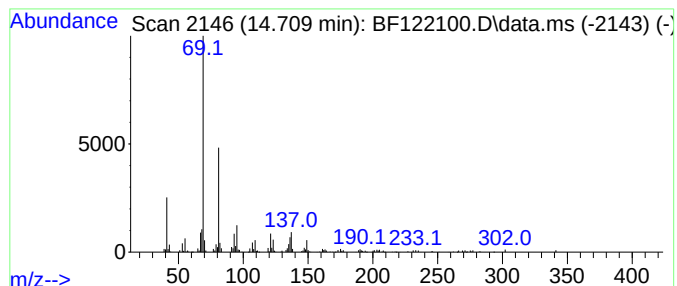
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Supraene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.710	5.76 ng	188275	Perylene-d12	15.327

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Supraene	410	C30H50	007683-64-9	91
2			Squalene	410	C30H50	000111-02-4	91
3			7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	83
4			5,9,13-Pentadecatrien-2-one, 6,1...	262	C18H30O	001117-52-8	64
5			2,3,5-Trimethyl-2,3,5-hexanetric...	203	C12H17N3	003974-79-6	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102320\
Data File : BF122100.D
Acq On : 23 Oct 2020 15:46
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Sample : L4488-05
Misc :
ALS Vial : 6 Sample Multiplier: 1

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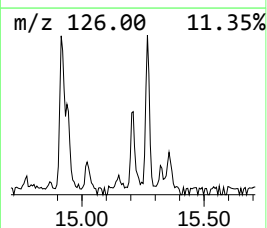
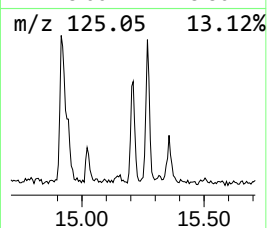
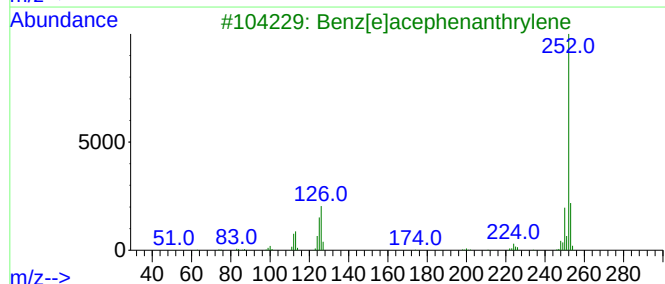
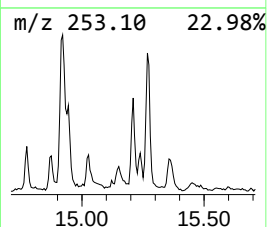
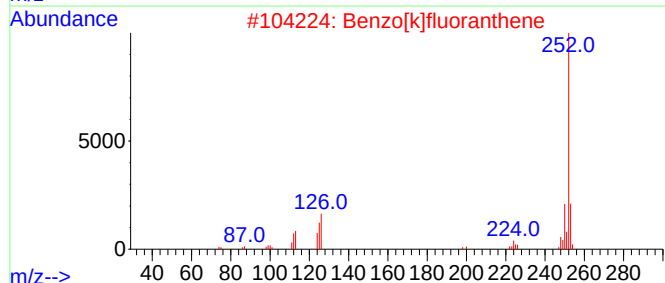
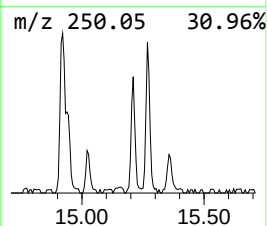
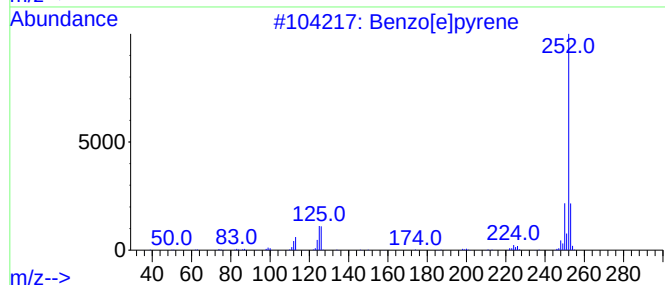
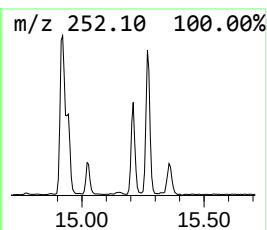
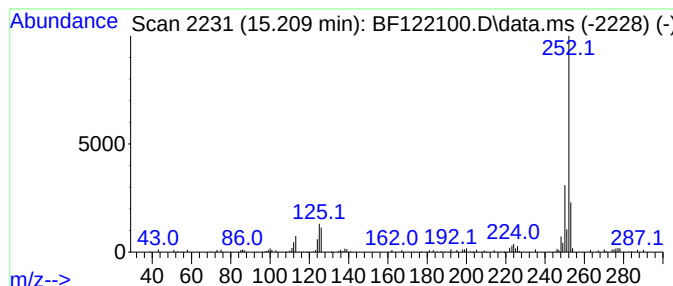
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzo[e]pyrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.210	2.89 ng	94371	Perylene-d12	15.327

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	98
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
3			Benz[e]acephenanthrylene	252	C20H12	000205-99-2	95
4			Perylene	252	C20H12	000198-55-0	95
5			Benzo[a]pyrene	252	C20H12	000050-32-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102320\
Data File : BF122100.D
Acq On : 23 Oct 2020 15:46
Operator : JU/CG
Sample : L4488-05
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.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
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Tetradecane	11.139	2.0	ng	100854	4	11.245	984634	20.0
1H-Benzotriazol...	11.492	2.1	ng	103420	4	11.245	984634	20.0
Phenanthrene, 1...	11.780	4.8	ng	238708	4	11.245	984634	20.0
Octadecanoic acid	12.563	8.6	ng	425303	4	11.245	984634	20.0
Phenanthrene, 4...	12.616	2.0	ng	93366	5	13.892	910658	20.0
11H-Benzo[a]flu...	13.751	5.8	ng	262112	5	13.892	910658	20.0
Supraene	14.710	5.8	ng	188275	6	15.327	653571	20.0
Benzo[e]pyrene	15.210	2.9	ng	94371	6	15.327	653571	20.0