

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF072720\
 Data File : BF121035.D
 Acq On : 27 Jul 2020 16:42
 Operator : JU/CG
 Sample : L3382-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CTR-007

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-BF071620.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	5.028	496	500	508	rVB	63581	79216	1.80%	0.192%
2	5.428	563	568	571	rBV	3071403	3031684	68.71%	7.344%
3	6.445	735	741	744	rBV	2852041	3059683	69.35%	7.412%
4	6.639	771	774	787	rBV	132575	158139	3.58%	0.383%
5	6.810	800	803	807	rBV	1381515	1163835	26.38%	2.819%
6	7.375	894	899	902	rBV	2223222	2133464	48.35%	5.168%
7	8.092	1017	1021	1025	rBV	1644445	1514428	34.32%	3.669%
8	9.169	1199	1204	1208	rBV	3526997	3884453	88.04%	9.410%
9	9.286	1222	1224	1233	rVB	42712	47838	1.08%	0.116%
10	9.563	1267	1271	1279	rBV	597595	534997	12.13%	1.296%
11	9.851	1314	1320	1323	rBV	1704470	1713706	38.84%	4.151%
12	10.633	1448	1453	1471	rBV	2229992	2817388	63.85%	6.825%
13	10.757	1471	1474	1482	rVB7	49125	82335	1.87%	0.199%
14	11.333	1567	1572	1574	rBV	2038600	1900496	43.07%	4.604%
15	11.357	1574	1576	1580	rVV	401825	356888	8.09%	0.865%
16	11.404	1582	1584	1588	rVB	76830	68536	1.55%	0.166%
17	11.663	1625	1628	1632	rVB	54691	47092	1.07%	0.114%
18	11.833	1653	1657	1659	rBV	152842	136493	3.09%	0.331%
19	11.863	1659	1662	1666	rVB	350684	319797	7.25%	0.775%
20	11.939	1672	1675	1678	rBV2	149703	162459	3.68%	0.394%
21	11.968	1678	1680	1685	rVB	96636	86027	1.95%	0.208%
22	12.127	1699	1707	1709	rBV3	77139	84366	1.91%	0.204%
23	12.151	1709	1711	1716	rVB2	52473	57891	1.31%	0.140%
24	12.310	1735	1738	1740	rBV	76384	69536	1.58%	0.168%
25	12.380	1744	1750	1753	rBV	185031	192576	4.36%	0.467%
26	12.416	1753	1756	1759	rVV2	91102	121394	2.75%	0.294%
27	12.463	1762	1764	1769	rBV3	81909	97289	2.21%	0.236%
28	12.539	1774	1777	1782	rVB	446817	416985	9.45%	1.010%
29	12.586	1782	1785	1787	rBV	49384	50922	1.15%	0.123%
30	12.645	1792	1795	1798	rBV3	44846	63348	1.44%	0.153%
31	12.768	1810	1816	1819	rBV	589808	552632	12.53%	1.339%
32	12.827	1823	1826	1830	rVB2	84018	97161	2.20%	0.235%
33	12.886	1830	1836	1837	rBV3	43454	59037	1.34%	0.143%
34	12.921	1837	1842	1845	rBV2	4017042	4412169	100.00%	10.688%

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If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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35	12.992	1850	1854	1861	rVB3	99166	166981	3.78%	0.405%
36	13.080	1865	1869	1871	rBV3	94156	141877	3.22%	0.344%
37	13.098	1871	1872	1876	rVB	115752	74728	1.69%	0.181%
38	13.163	1878	1883	1886	rBV3	78827	111483	2.53%	0.270%
39	13.204	1886	1890	1898	rVB2	155158	161791	3.67%	0.392%
40	13.298	1902	1906	1908	rBV	134821	122716	2.78%	0.297%
41	13.327	1908	1911	1914	rVB	93593	87256	1.98%	0.211%
42	13.398	1920	1923	1929	rVB4	82277	100332	2.27%	0.243%
43	13.492	1937	1939	1947	rVB7	40070	69374	1.57%	0.168%
44	13.604	1956	1958	1963	rVB	96635	104731	2.37%	0.254%
45	13.715	1972	1977	1980	rBV3	95127	158358	3.59%	0.384%
46	13.768	1984	1986	1991	rVV2	60805	83371	1.89%	0.202%
47	13.833	1991	1997	2012	rVB4	326976	853120	19.34%	2.067%
48	13.968	2014	2020	2023	rBV	2026994	2307109	52.29%	5.589%
49	13.992	2023	2024	2032	rVB	488295	369049	8.36%	0.894%
50	14.133	2045	2048	2054	rVB5	64958	83039	1.88%	0.201%
51	14.221	2061	2063	2066	rVB2	57477	46471	1.05%	0.113%
52	14.351	2081	2085	2089	rVV	152191	223038	5.06%	0.540%
53	14.386	2089	2091	2095	rVV2	126001	136762	3.10%	0.331%
54	14.439	2095	2100	2103	rVV3	125882	199830	4.53%	0.484%
55	14.480	2103	2107	2115	rVV3	129532	280021	6.35%	0.678%
56	14.551	2117	2119	2124	rVB2	51430	56615	1.28%	0.137%
57	14.686	2139	2142	2147	rVB2	30161	53986	1.22%	0.131%
58	14.739	2148	2151	2154	rBV2	96175	112321	2.55%	0.272%
59	14.815	2159	2164	2175	rVV	570816	835086	18.93%	2.023%
60	14.904	2177	2179	2184	rVB2	52700	57396	1.30%	0.139%
61	14.998	2190	2195	2198	rBV	503650	767566	17.40%	1.859%
62	15.074	2205	2208	2211	rBV	122716	122785	2.78%	0.297%
63	15.204	2224	2230	2234	rBV4	43371	96529	2.19%	0.234%
64	15.292	2241	2245	2251	rVB	328960	398832	9.04%	0.966%
65	15.351	2251	2255	2261	rVB	224800	296729	6.73%	0.719%
66	15.415	2261	2266	2270	rBV	1204668	1607386	36.43%	3.894%
67	15.445	2270	2271	2279	rVB2	186360	215777	4.89%	0.523%
68	15.762	2323	2325	2332	rVB3	36046	50346	1.14%	0.122%
69	15.874	2340	2344	2350	rVB	102036	143861	3.26%	0.349%
70	15.968	2356	2360	2362	rBV4	30537	51437	1.17%	0.125%
71	16.368	2423	2428	2432	rBV3	45254	74748	1.69%	0.181%

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

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72	16.662	2471	2478	2480	rBV3	32967	71418	1.62%	0.173%
73	16.839	2502	2508	2518	rVB2	125598	265008	6.01%	0.642%
74	16.939	2522	2525	2532	rVB4	36193	71476	1.62%	0.173%
75	17.015	2533	2538	2542	rBV2	29611	56272	1.28%	0.136%
76	17.068	2543	2547	2551	rBV3	32944	60120	1.36%	0.146%
77	17.186	2563	2567	2573	rVB7	42619	79281	1.80%	0.192%
78	17.262	2573	2580	2592	rVB	96931	223143	5.06%	0.541%
79	17.621	2638	2641	2650	rVB6	24834	56241	1.27%	0.136%

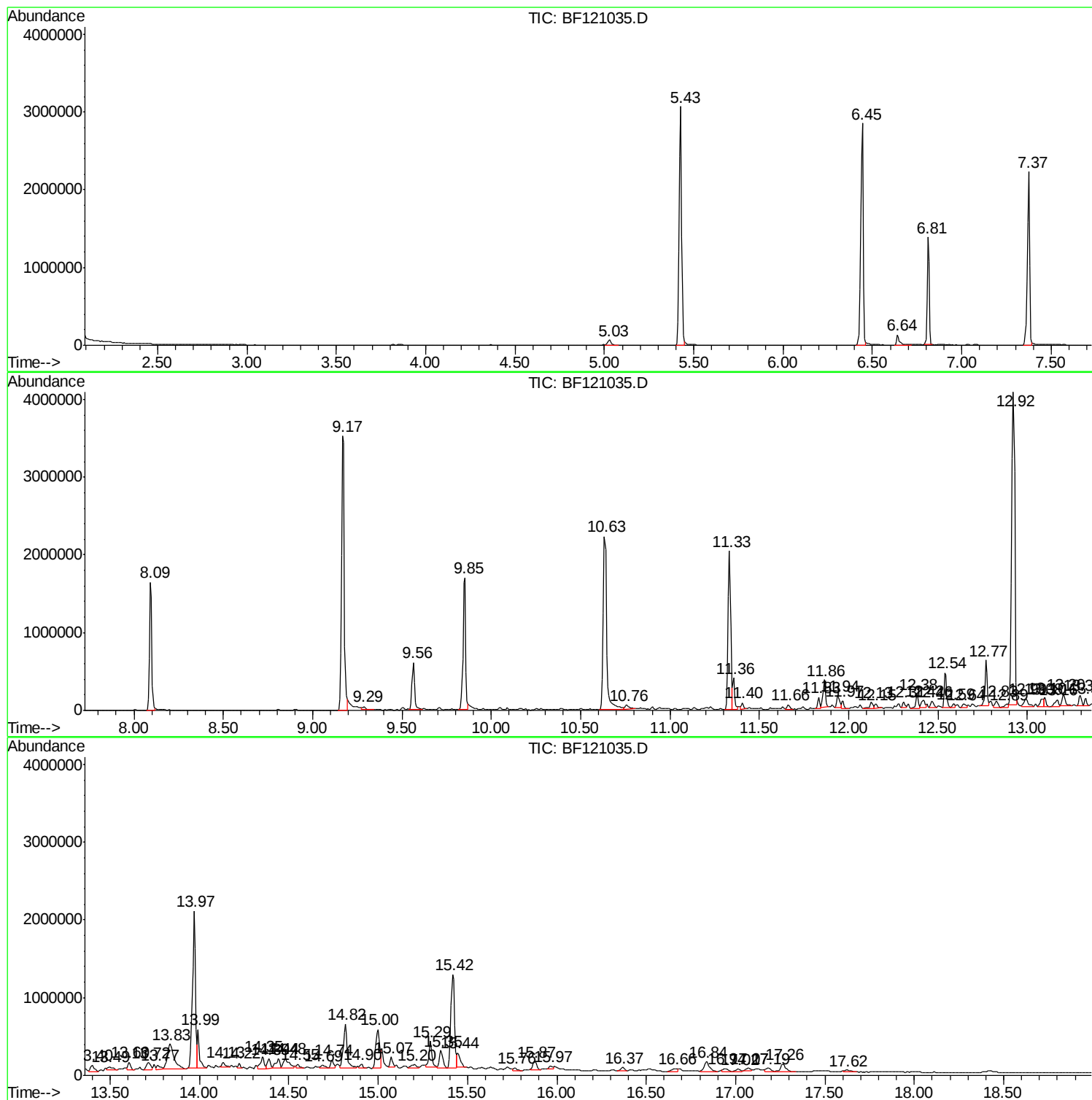
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Instrument :
BNA_F
ClientSampled :
CTR-007

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF071620.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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Operator : JU/CG
Sample : L3382-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CTR-007

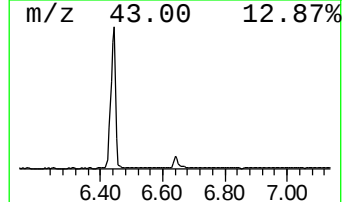
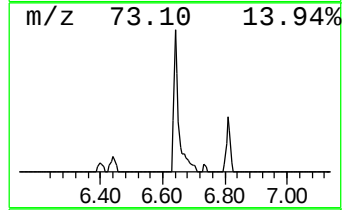
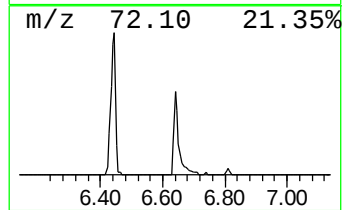
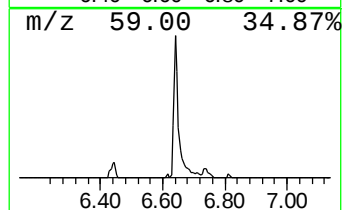
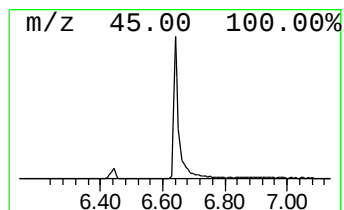
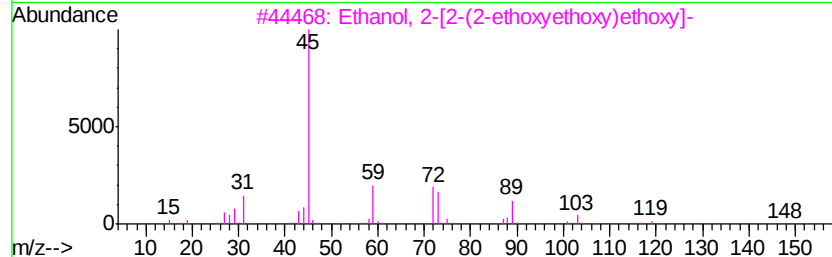
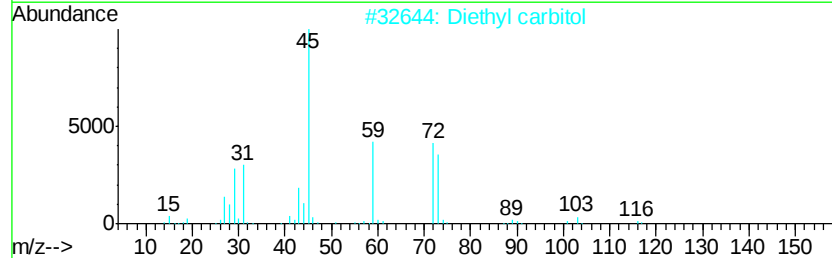
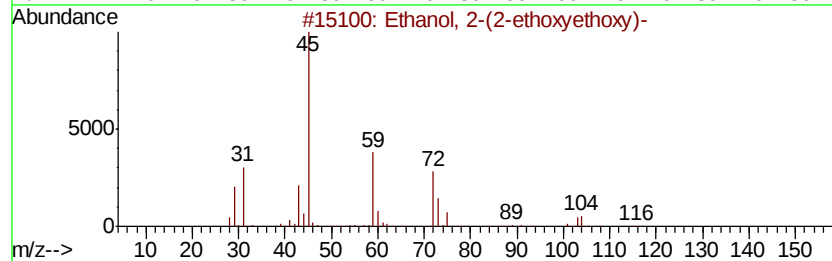
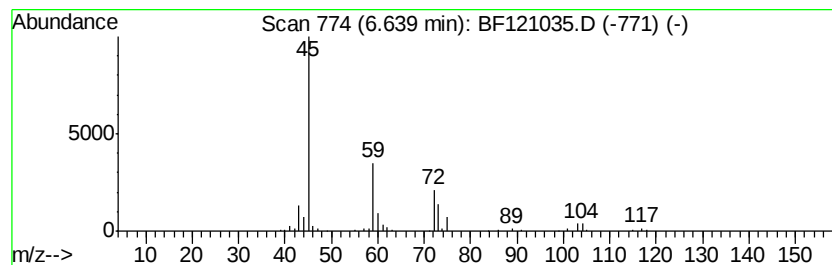
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.64	2.72 ng	158139	1,4-Dichlorobenzene-d4	6.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	90
2		Diethyl carbitol	162	C8H18O3	000112-36-7	72
3		Ethanol, 2-[2-(2-ethoxyethoxy)et...	178	C8H18O4	000112-50-5	64
4		2-Propanol, 1-(1-methylethoxy)-	118	C6H14O2	003944-36-3	45
5		Ethanol, 2-(2-methoxyethoxy)-	120	C5H12O3	000111-77-3	42



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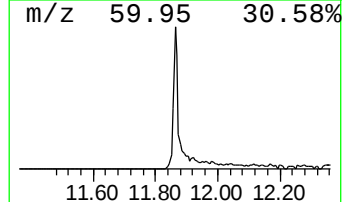
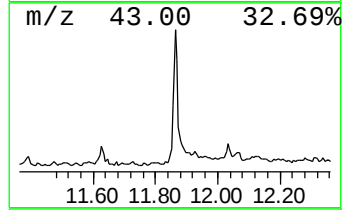
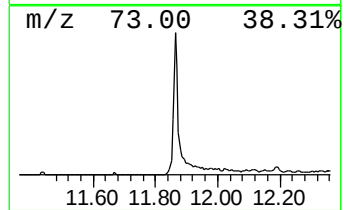
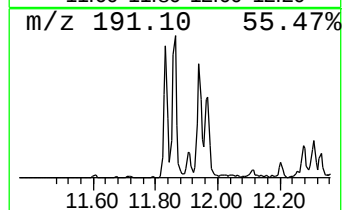
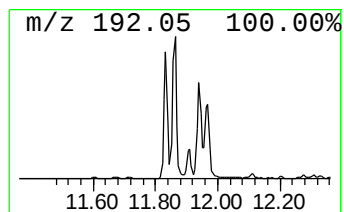
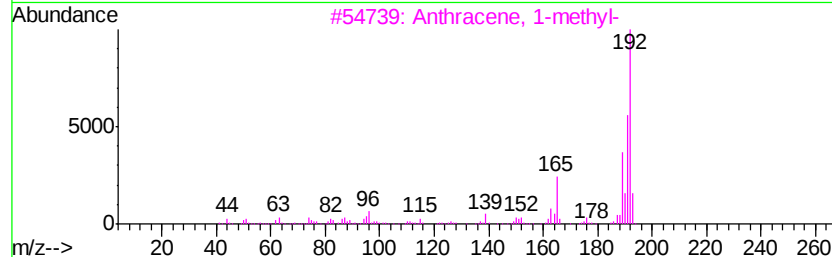
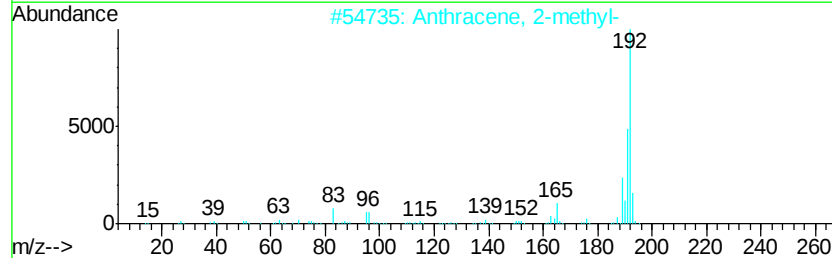
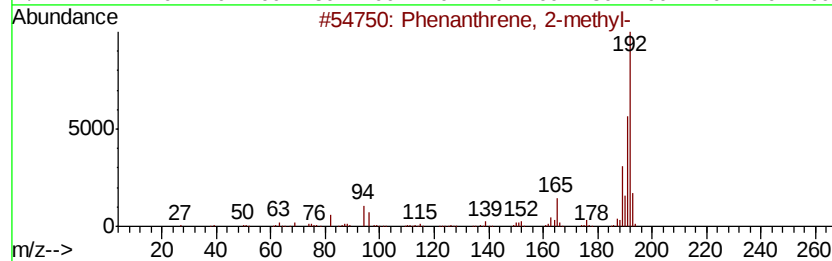
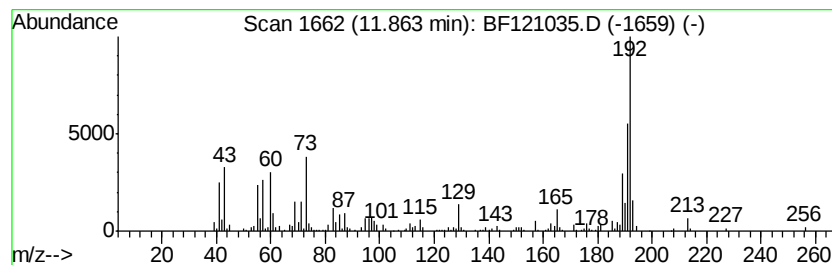
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Phenanthrene, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.86	3.37 ng	319797	Phenanthrene-d10	11.33	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	96
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	95



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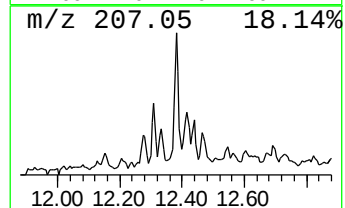
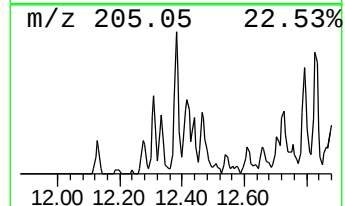
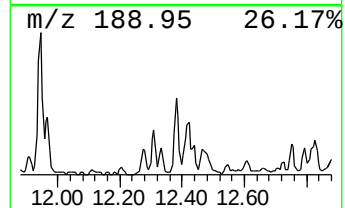
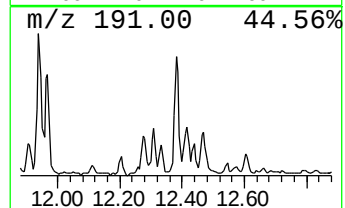
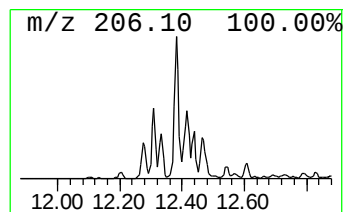
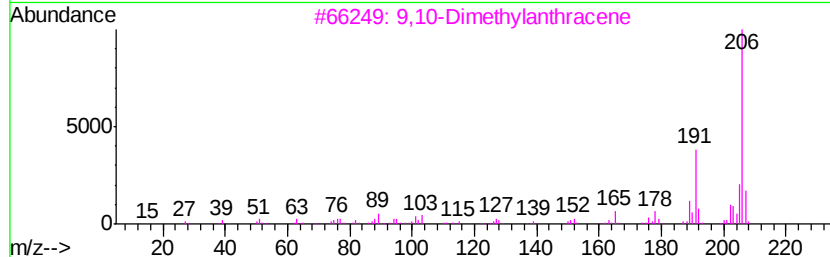
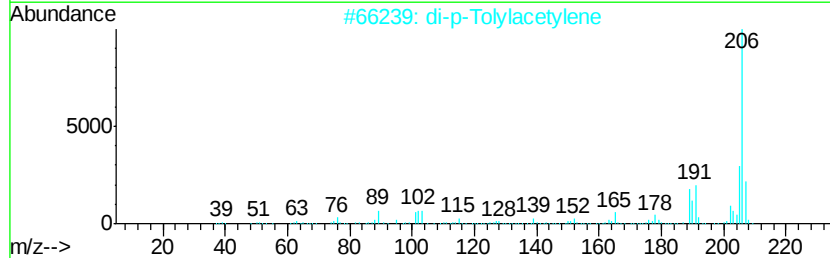
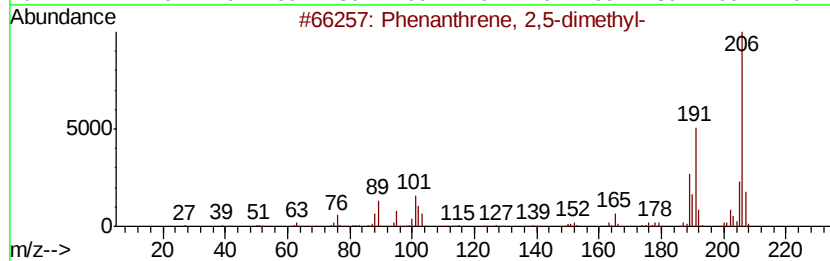
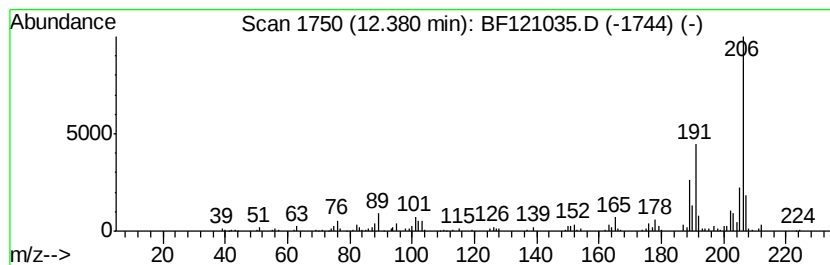
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Phenanthrene, 2,5-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.38	2.03 ng	192576	Phenanthrene-d10		11.33
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	97
2	di-p-Tolylacetylvene	206	C16H14	002789-88-0	96
3	9,10-Dimethylanthracene	206	C16H14	000781-43-1	96
4	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	93
5	Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	93



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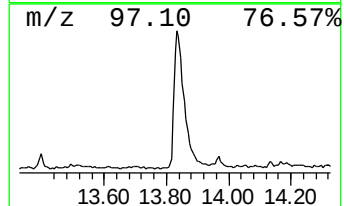
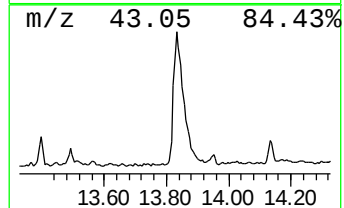
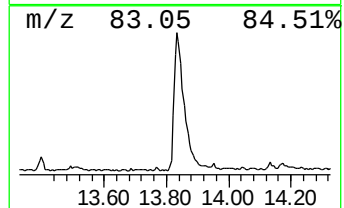
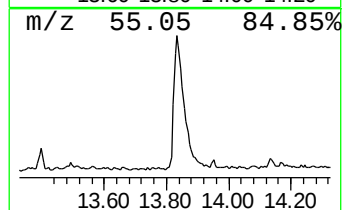
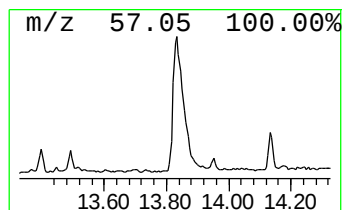
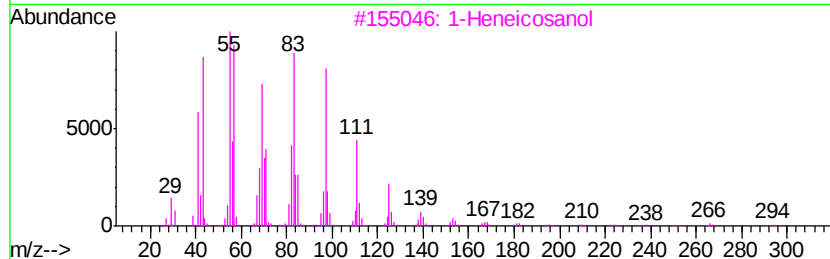
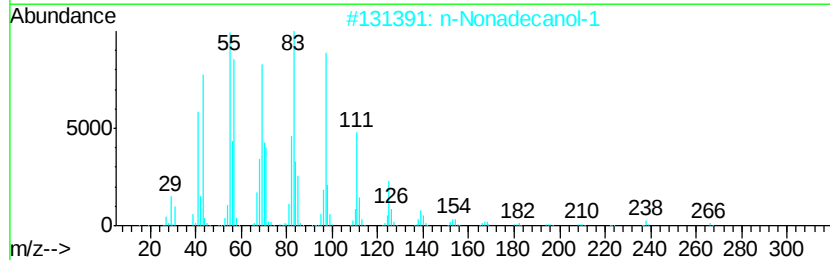
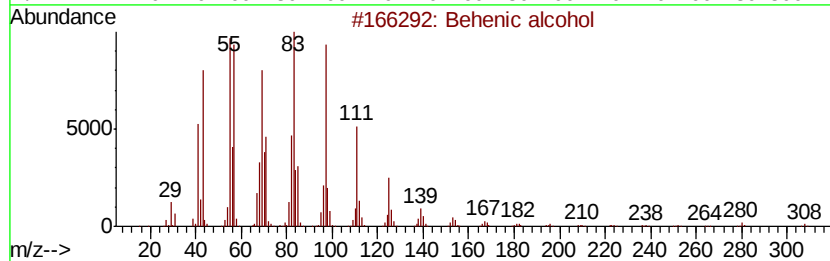
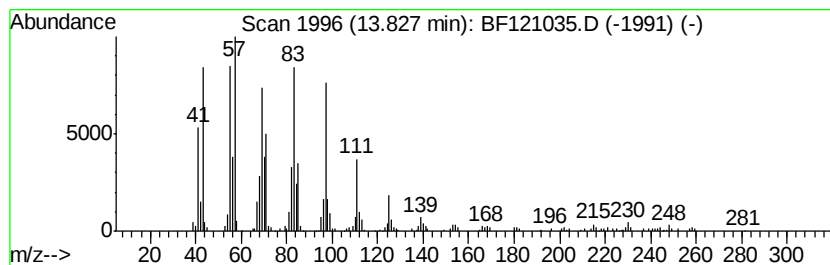
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Behenic alcohol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.83	7.40 ng	853120	Chrysene-d12	13.97

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Behenic alcohol	326	C22H46O	000661-19-8	95
2		n-Nonadecanol-1	284	C19H40O	001454-84-8	95
3		1-Heneicosanol	312	C21H44O	015594-90-8	95
4		n-Tetracosanol-1	354	C24H50O	000506-51-4	95
5		1-Octadecene	252	C18H36	000112-88-9	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072720\
Data File : BF121035.D
Acq On : 27 Jul 2020 16:42
Operator : JU/CG
Sample : L3382-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CTR-007

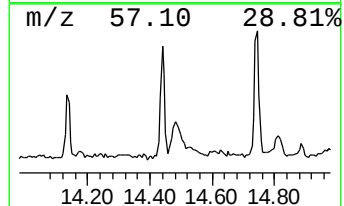
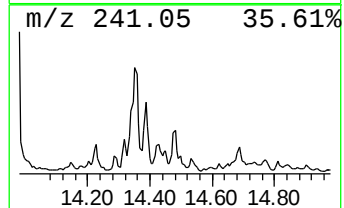
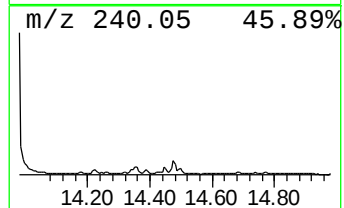
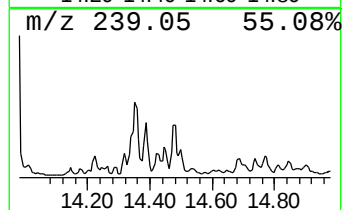
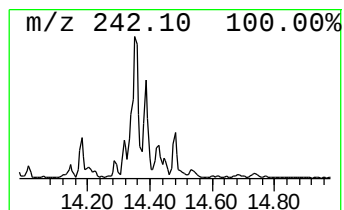
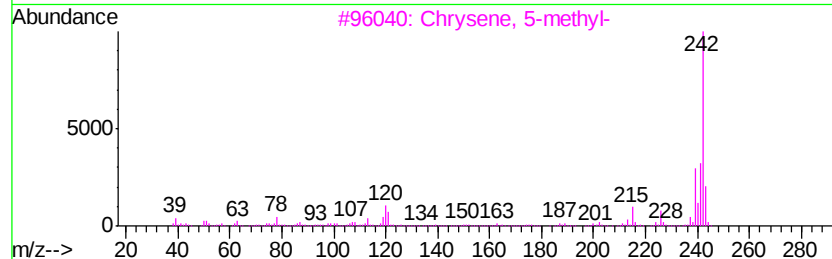
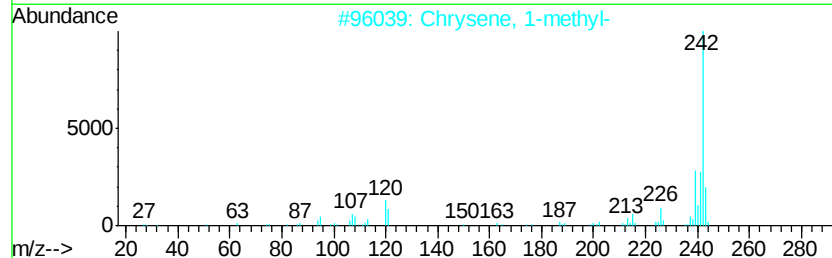
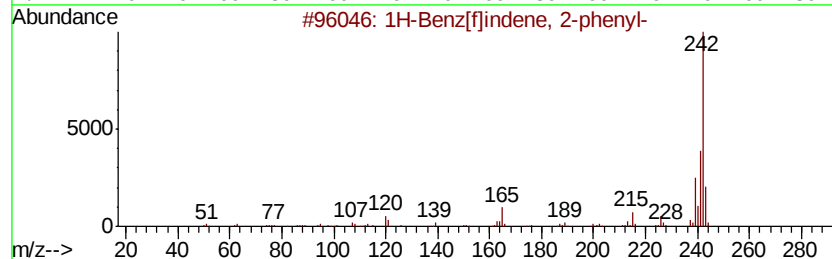
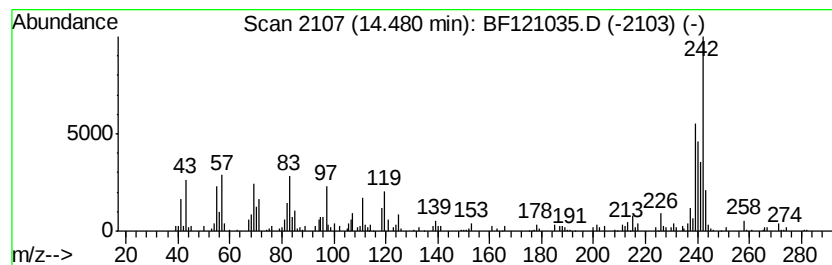
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 1H-Benz[f]indene, 2-phenyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.48	2.43 ng	280021	Chrysene-d12	13.97

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Benz[f]indene, 2-phenyl-	242	C19H14	1000305-22-4	89
2		Chrysene, 1-methyl-	242	C19H14	003351-28-8	64
3		Chrysene, 5-methyl-	242	C19H14	003697-24-3	64
4		Benz[a]anthracene, 1-methyl-	242	C19H14	002498-77-3	64
5		Benz[a]anthracene, 12-methyl-	242	C19H14	002422-79-9	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072720\
Data File : BF121035.D
Acq On : 27 Jul 2020 16:42
Operator : JU/CG
Sample : L3382-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CTR-007

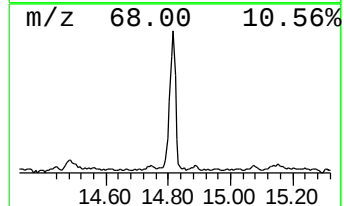
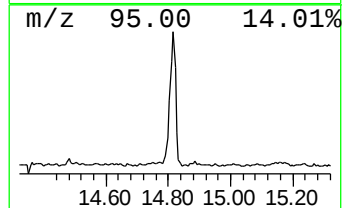
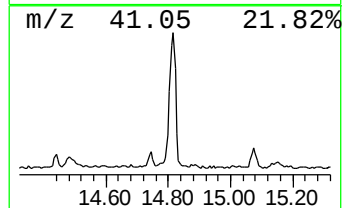
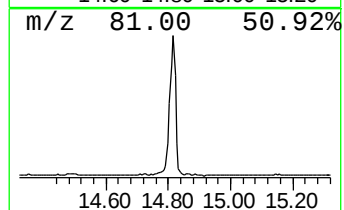
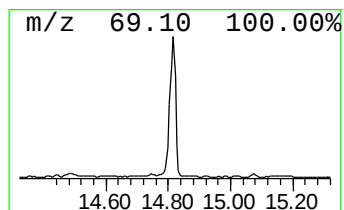
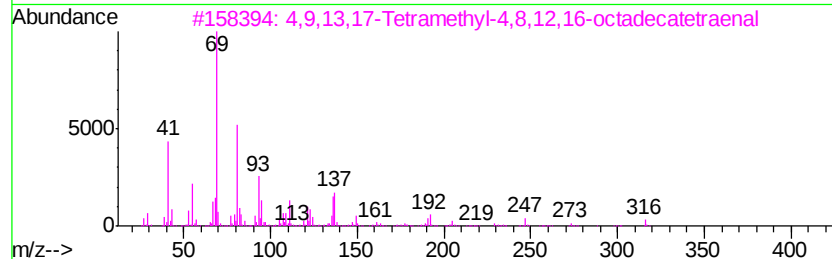
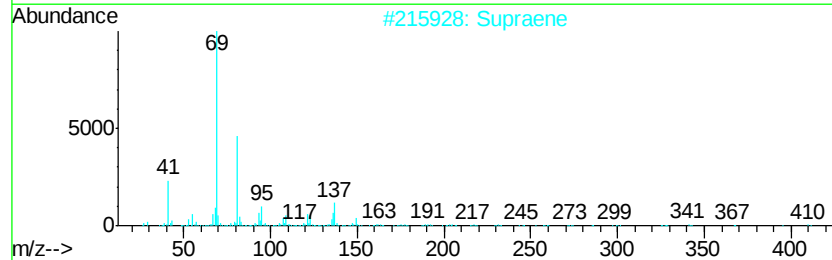
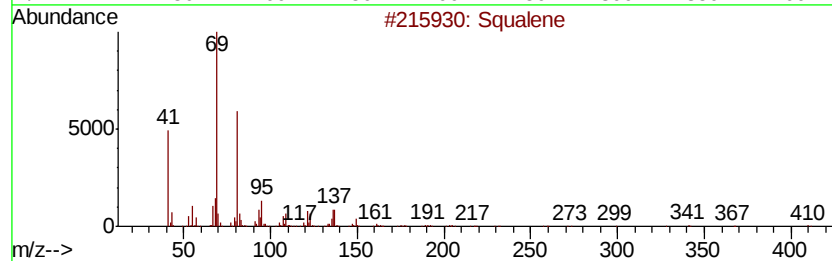
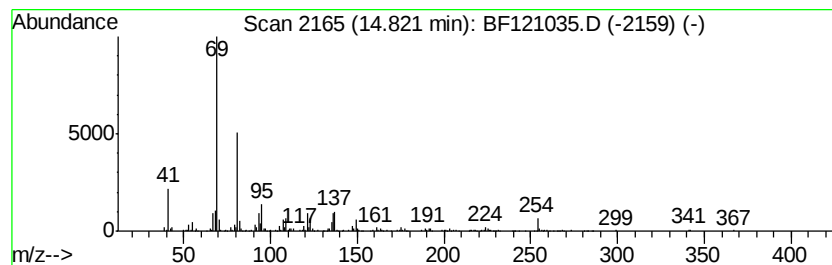
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Squalene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	10.39 ng	835086	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	000111-02-4	91
2		Supraene	410	C30H50	007683-64-9	91
3		4,9,13,17-Tetramethyl-4,8,12,16-...	316	C22H36O	056882-09-8	83
4		7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	72
5		2,3,5-Trimethyl-2,3,5-hexanetric...	203	C12H17N3	003974-79-6	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072720\
Data File : BF121035.D
Acq On : 27 Jul 2020 16:42
Operator : JU/CG
Sample : L3382-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

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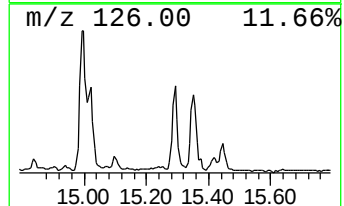
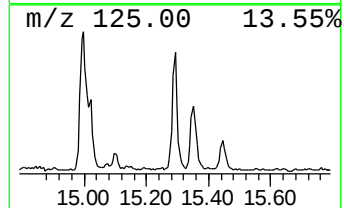
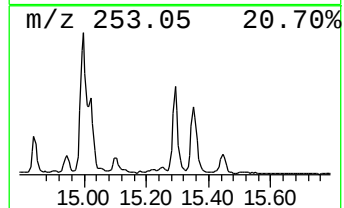
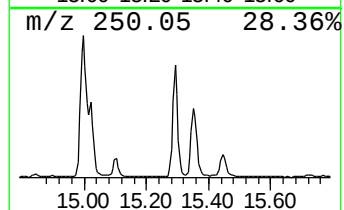
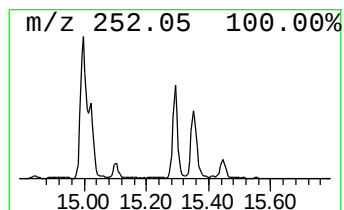
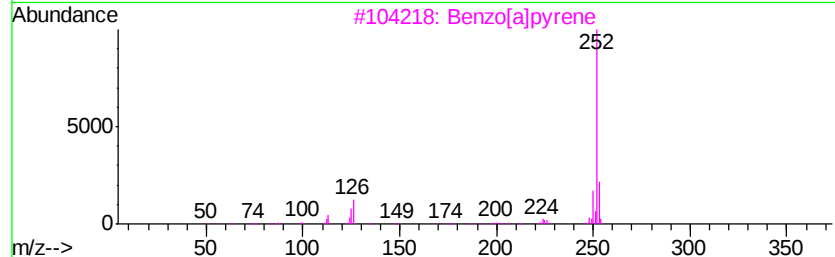
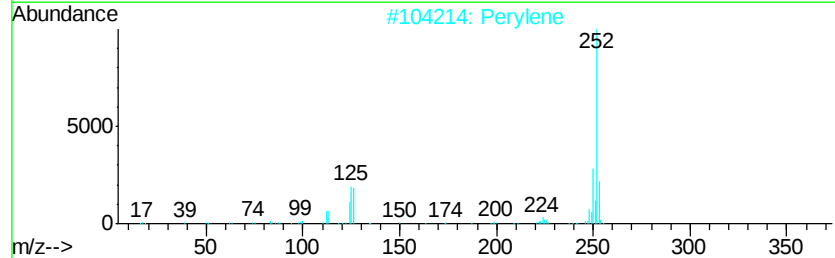
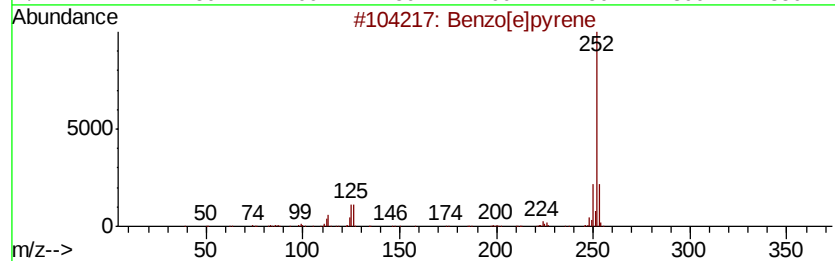
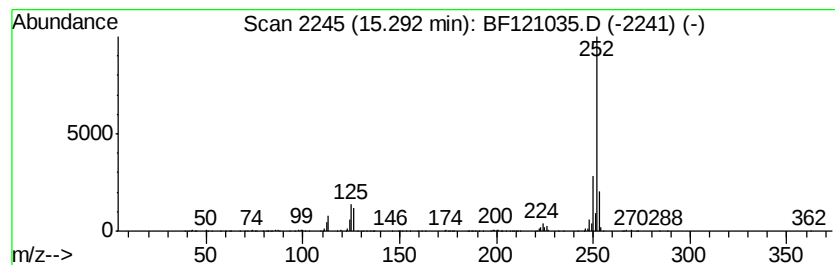
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.29	4.96 ng	398832	Perylene-d12	15.42

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	99
2			Perylene	252	C20H12	000198-55-0	95
3			Benzo[a]pyrene	252	C20H12	000050-32-8	94
4			Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	93
5			Benzo[k]fluoranthene	252	C20H12	000207-08-9	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF072720\
Data File : BF121035.D
Acq On : 27 Jul 2020 16:42
Operator : JU/CG
Sample : L3382-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CTR-007

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF071620.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
Ethanol, 2-(2-eth...	6.64	2.7	ng	158139	1	6.81	1163840	20.0
Phenanthrene, 2-m...	11.86	3.4	ng	319797	4	11.33	1900500	20.0
Phenanthrene, 2,5...	12.38	2.0	ng	192576	4	11.33	1900500	20.0
Behenic alcohol	13.83	7.4	ng	853120	5	13.97	2307110	20.0
1H-Benz[f]indene,...	14.48	2.4	ng	280021	5	13.97	2307110	20.0
Squalene	14.82	10.4	ng	835086	6	15.42	1607390	20.0
Benzo[e]pyrene	15.29	5.0	ng	398832	6	15.42	1607390	20.0