

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\
 Data File : BF121105.D
 Acq On : 29 Jul 2020 16:52
 Operator : JU/CG
 Sample : L3430-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS2

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	5.381	556	560	562	rBV	73720	82968	1.24%	0.111%
2	5.416	562	566	570	rVV	2719962	2863523	42.93%	3.815%
3	6.440	735	740	743	rBV	2783120	2900900	43.49%	3.865%
4	6.634	769	773	777	rBV2	63777	68978	1.03%	0.092%
5	6.798	798	801	805	rBV	1203625	984318	14.76%	1.312%
6	7.210	866	871	882	rBV4	39656	69015	1.03%	0.092%
7	7.363	890	897	900	rBV	2126387	2040947	30.60%	2.719%
8	8.081	1015	1019	1022	rBV	1413760	1253079	18.79%	1.670%
9	9.157	1198	1202	1206	rBV	3565881	3453150	51.77%	4.601%
10	9.298	1224	1226	1232	rVB	89300	95684	1.43%	0.127%
11	9.428	1243	1248	1252	rBV5	63002	91762	1.38%	0.122%
12	9.492	1257	1259	1262	rVV	165366	146174	2.19%	0.195%
13	9.551	1266	1269	1273	rVB	624462	489181	7.33%	0.652%
14	9.698	1289	1294	1299	rBV2	237577	395978	5.94%	0.528%
15	9.833	1312	1317	1320	rBV	1587154	1480146	22.19%	1.972%
16	9.863	1320	1322	1326	rVB2	179956	169611	2.54%	0.226%
17	9.957	1336	1338	1341	rVV2	114894	103596	1.55%	0.138%
18	9.986	1341	1343	1346	rVB2	90377	92101	1.38%	0.123%
19	10.039	1348	1352	1355	rBV	153600	137935	2.07%	0.184%
20	10.239	1383	1386	1388	rBV3	60218	74920	1.12%	0.100%
21	10.380	1407	1410	1414	rBV	150952	138408	2.08%	0.184%
22	10.539	1434	1437	1439	rBV3	82679	96736	1.45%	0.129%
23	10.592	1443	1446	1447	rBV	143108	119669	1.79%	0.159%
24	10.622	1447	1451	1455	rVB	1984486	1971381	29.56%	2.627%
25	10.745	1468	1472	1476	rBV4	169167	220491	3.31%	0.294%
26	10.869	1490	1493	1499	rBV5	89576	133825	2.01%	0.178%
27	10.928	1500	1503	1505	rBV3	81451	86269	1.29%	0.115%
28	10.986	1510	1513	1516	rBV4	138385	153998	2.31%	0.205%
29	11.051	1521	1524	1528	rBV4	67976	115032	1.72%	0.153%
30	11.122	1533	1536	1540	rVB	147601	136061	2.04%	0.181%
31	11.186	1541	1547	1550	rBV5	97379	162740	2.44%	0.217%
32	11.216	1550	1552	1555	rVV	85588	76709	1.15%	0.102%
33	11.275	1559	1562	1565	rBV3	72409	106439	1.60%	0.142%
34	11.322	1566	1570	1572	rBV	1353598	1360944	20.40%	1.813%

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35	11.345	1572	1574	1577	rVV	914277	819770	12.29%	1.092%
36	11.375	1577	1579	1580	rVV	186724	147837	2.22%	0.197%
37	11.392	1580	1582	1585	rVB	452846	358751	5.38%	0.478%
38	11.427	1585	1588	1591	rBV2	143160	159823	2.40%	0.213%
39	11.457	1591	1593	1597	rVB	184298	167062	2.50%	0.223%
40	11.539	1604	1607	1611	rBV2	146949	230953	3.46%	0.308%
41	11.616	1617	1620	1624	rBV	106086	89434	1.34%	0.119%
42	11.698	1631	1634	1638	rBV5	83516	130444	1.96%	0.174%
43	11.816	1650	1654	1657	rBV3	304203	373700	5.60%	0.498%
44	11.863	1657	1662	1665	rBV2	1479854	1692606	25.38%	2.255%
45	11.933	1671	1674	1676	rBV2	265170	253561	3.80%	0.338%
46	12.016	1683	1688	1690	rBV3	175950	294049	4.41%	0.392%
47	12.063	1694	1696	1699	rVB2	184324	153453	2.30%	0.204%
48	12.116	1700	1705	1707	rBV2	241762	332166	4.98%	0.443%
49	12.139	1707	1709	1713	rVB3	150676	170888	2.56%	0.228%
50	12.174	1713	1715	1717	rBV	89992	91188	1.37%	0.121%
51	12.263	1725	1730	1733	rBV2	291446	380051	5.70%	0.506%
52	12.298	1733	1736	1738	rBV	232488	207483	3.11%	0.276%
53	12.351	1742	1745	1746	rBV	138829	134597	2.02%	0.179%
54	12.369	1746	1748	1753	rBV3	166291	230886	3.46%	0.308%
55	12.457	1760	1763	1764	rBV	312803	312723	4.69%	0.417%
56	12.533	1772	1776	1779	rBV	2314987	2228925	33.42%	2.970%
57	12.574	1779	1783	1790	rVB	1872756	2165622	32.47%	2.885%
58	12.639	1791	1794	1799	rVV2	606693	673037	10.09%	0.897%
59	12.763	1809	1815	1818	rBV2	2077470	2292497	34.37%	3.055%
60	12.792	1818	1820	1830	rVB2	817888	1078224	16.17%	1.437%
61	12.910	1836	1840	1846	rVB	2723436	2881057	43.19%	3.839%
62	13.010	1855	1857	1860	rBV3	114641	91505	1.37%	0.122%
63	13.045	1860	1863	1865	rBV	495685	456901	6.85%	0.609%
64	13.086	1868	1870	1873	rVB	459368	449421	6.74%	0.599%
65	13.157	1879	1882	1884	rVB	322853	266502	4.00%	0.355%
66	13.192	1885	1888	1890	rVB2	281460	258586	3.88%	0.345%
67	13.280	1901	1903	1906	rBV2	219546	242190	3.63%	0.323%
68	13.410	1923	1925	1927	rVB	232287	165243	2.48%	0.220%
69	13.598	1954	1957	1961	rVB	226558	257932	3.87%	0.344%
70	13.739	1978	1981	1983	rBV	268758	263977	3.96%	0.352%
71	13.810	1990	1993	1998	rVB2	838453	899153	13.48%	1.198%

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 Stop Thrs : 0 Peak Location: TOP

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72	13.957	2013	2018	2021	rBV3	1735771	2756524	41.33%	3.673%
73	13.986	2021	2023	2028	rVB	1389153	1203784	18.05%	1.604%
74	14.445	2098	2101	2104	rVB2	538938	508636	7.63%	0.678%
75	14.998	2190	2195	2197	rBV	1194999	1829738	27.43%	2.438%
76	15.068	2203	2207	2210	rBV	949524	1080576	16.20%	1.440%
77	15.098	2210	2212	2217	rVB	652062	661506	9.92%	0.881%
78	15.292	2241	2245	2251	rVB	709519	969593	14.54%	1.292%
79	15.351	2251	2255	2261	rVB	1107377	1394391	20.91%	1.858%
80	15.415	2262	2266	2268	rBV	799800	1022313	15.33%	1.362%
81	15.862	2338	2342	2347	rVB	742634	1066261	15.99%	1.421%
82	16.845	2504	2509	2512	rBV2	412942	813958	12.20%	1.085%
83	17.015	2534	2538	2543	rBV3	358489	772986	11.59%	1.030%
84	17.445	2602	2611	2622	rBV5	1689400	6669997	100.00%	8.887%
85	17.886	2680	2686	2695	rBV3	597310	1951843	29.26%	2.601%
86	17.980	2698	2702	2712	rVB4	876667	2422358	36.32%	3.228%
87	18.221	2736	2743	2760	rBV3	983895	4525016	67.84%	6.029%
88	18.427	2773	2778	2790	rVB3	783478	2127901	31.90%	2.835%

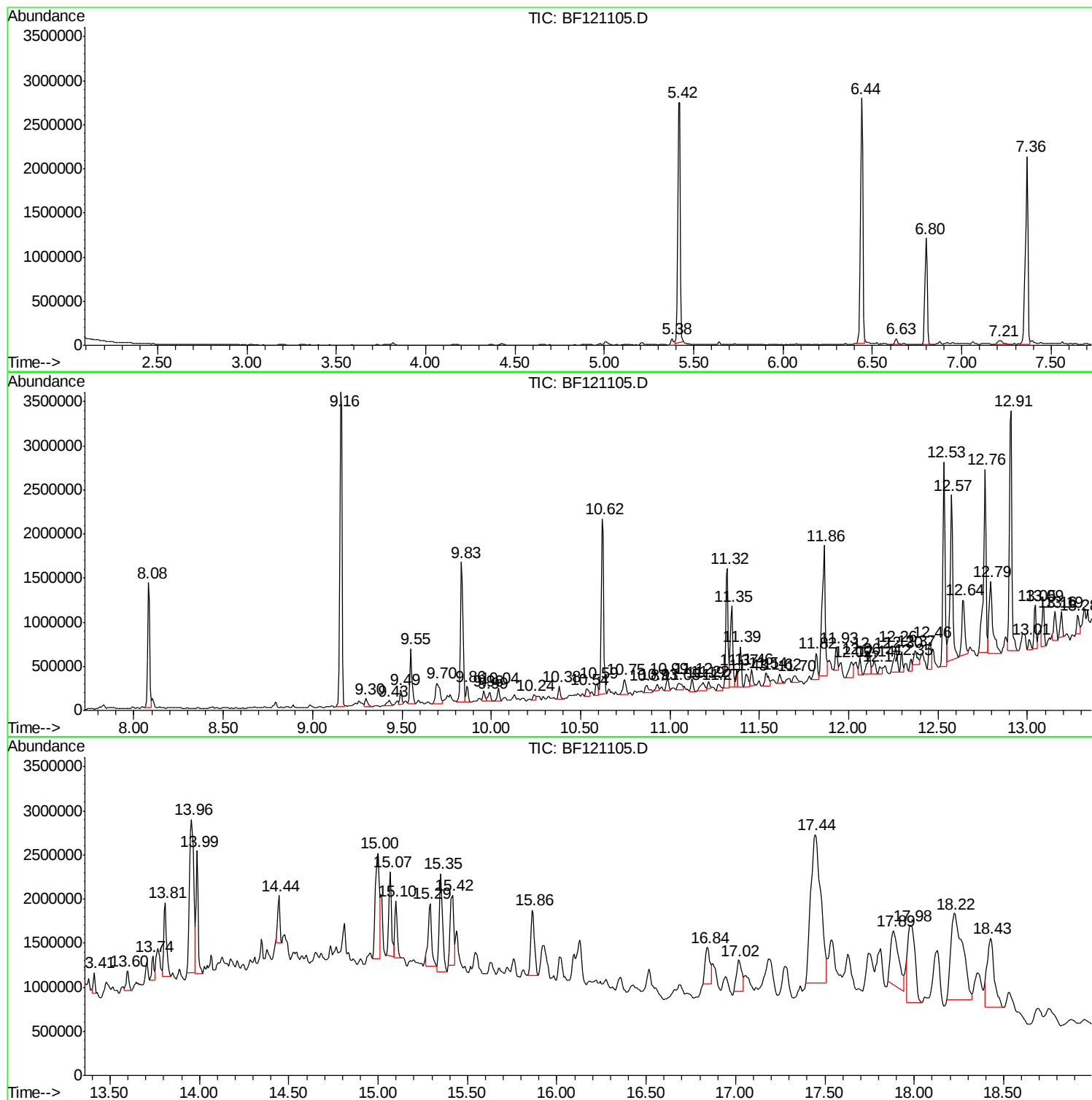
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ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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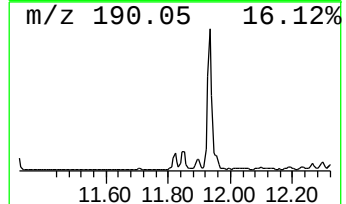
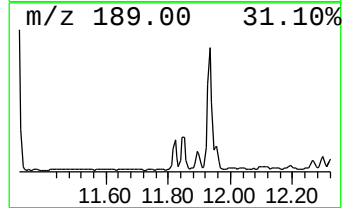
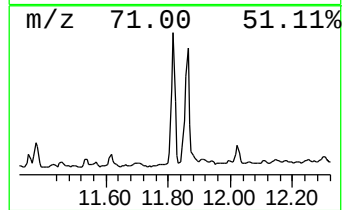
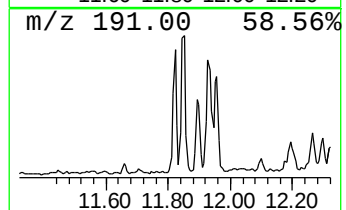
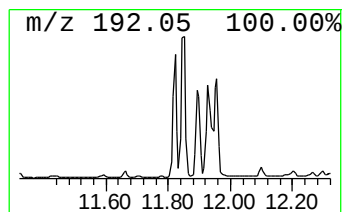
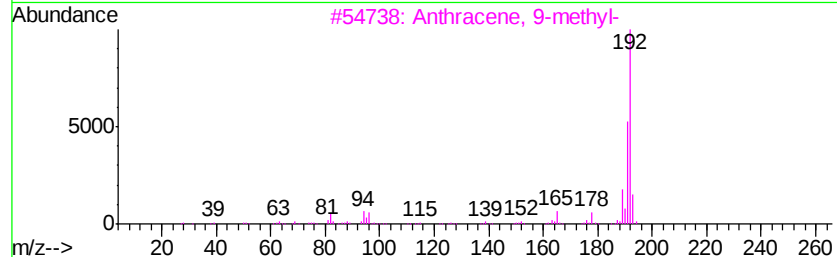
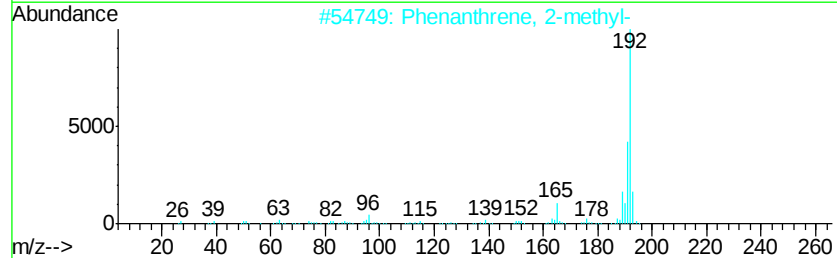
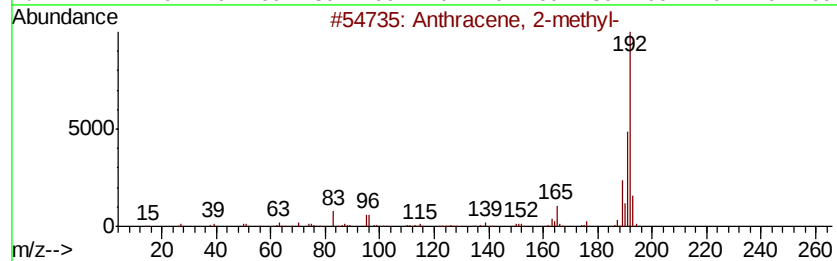
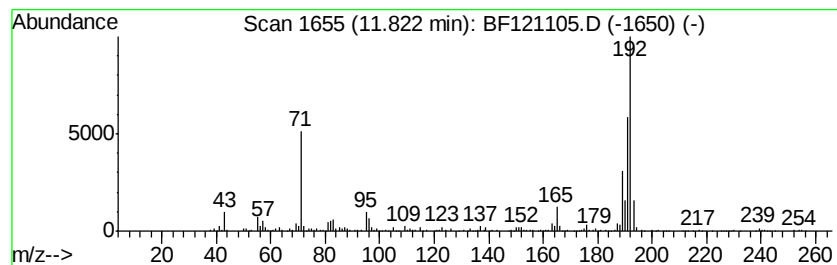
Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Anthracene, 2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.82	5.49 ng	373700	Phenanthrene-d10		11.32
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94
3	Anthracene, 9-methyl-	192	C15H12	000779-02-2	86
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	83
5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	52



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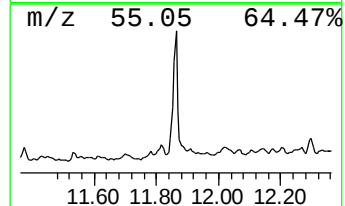
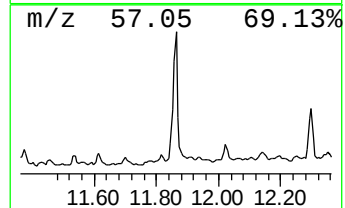
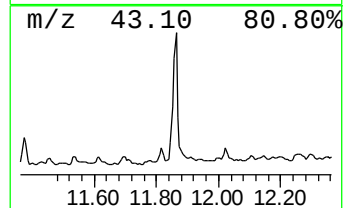
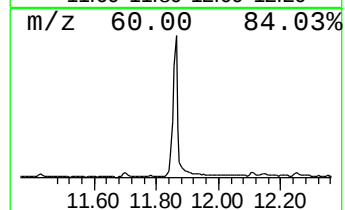
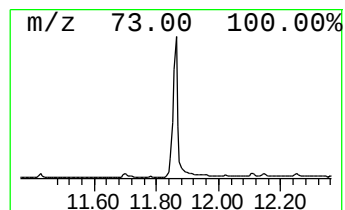
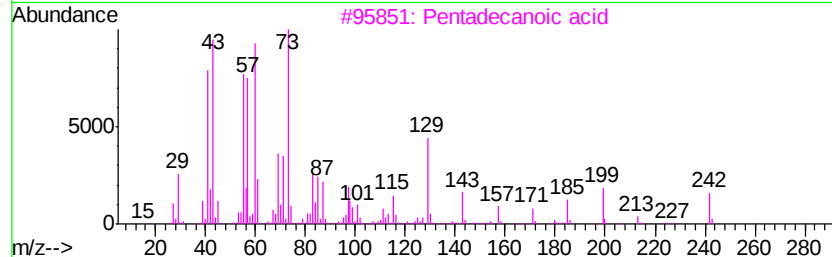
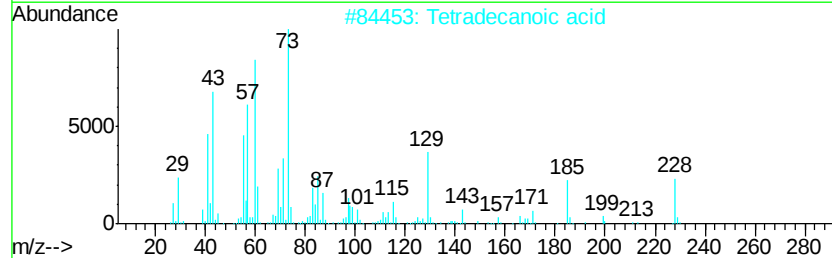
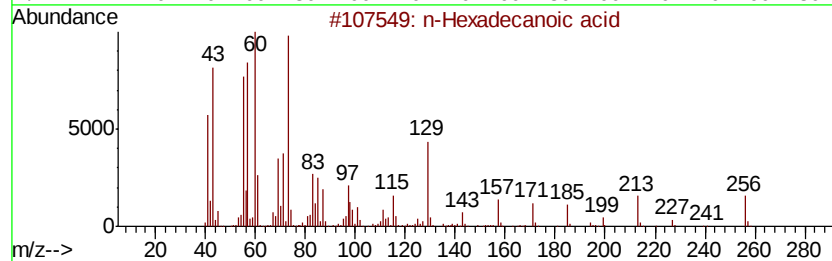
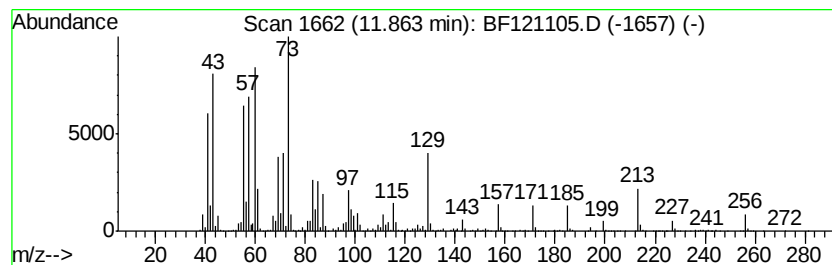
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.86	24.87 ng	1692610	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	94
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	93
4		Tridecanoic acid	214	C13H26O2	000638-53-9	74
5		n-Decanoic acid	172	C10H20O2	000334-48-5	68



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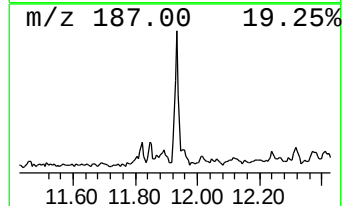
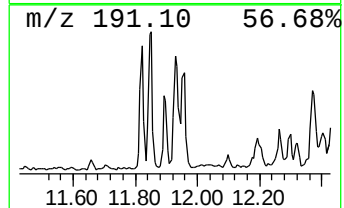
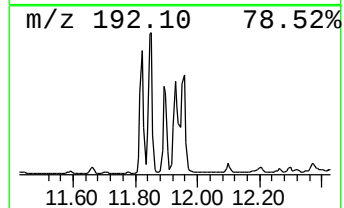
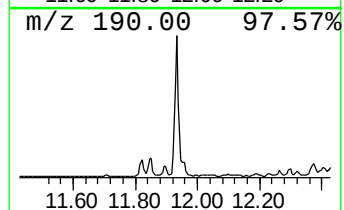
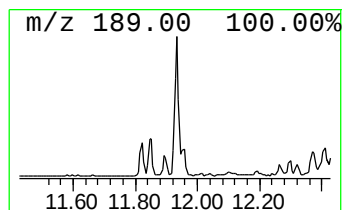
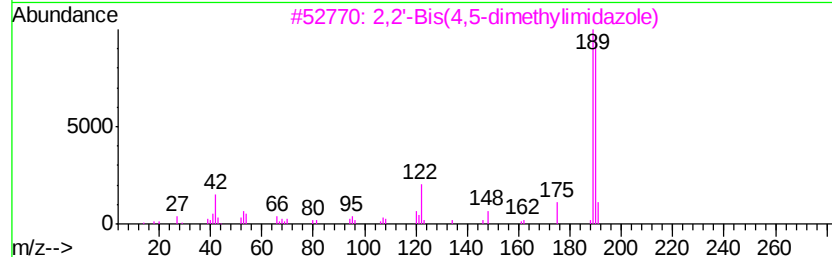
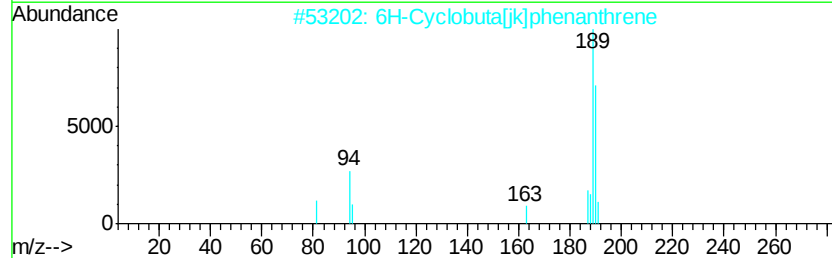
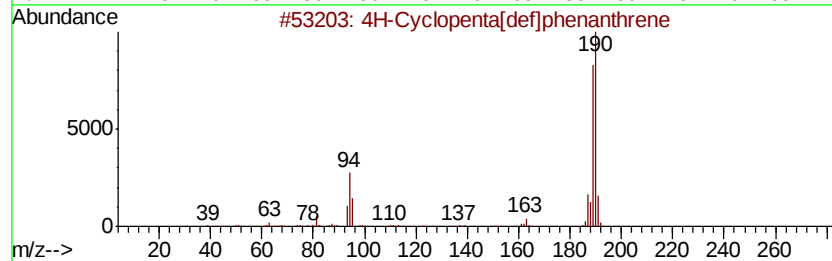
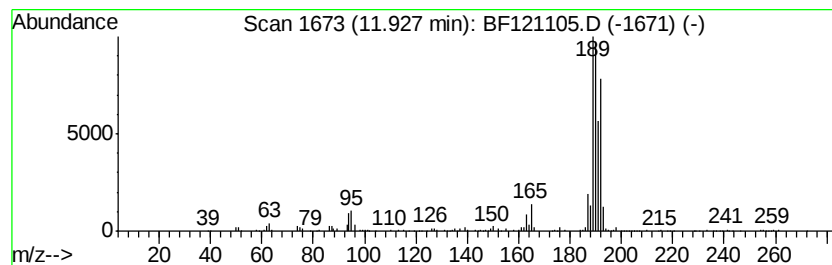
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 4H-Cyclopenta[def]phenanthrene Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	3.73 ng	253561	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	81
2		6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	64
3		2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	42
4		1,4-Naphthalenedione, 2,5-dihydr...	190	C10H6O4	004923-55-1	25
5		1-(Phenylethynyl)-1H-1,2,3-benzo...	219	C14H9N3	209912-23-2	17



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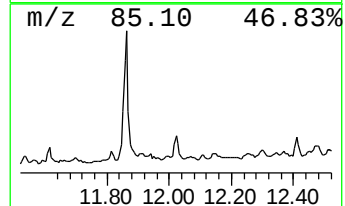
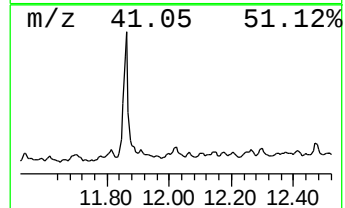
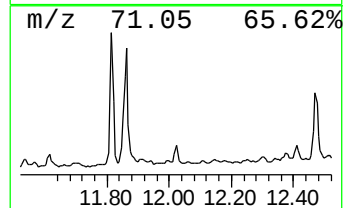
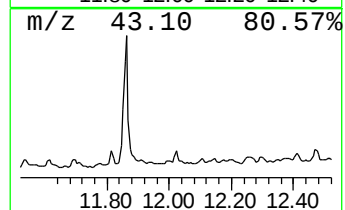
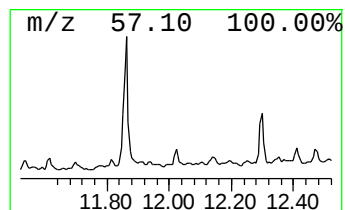
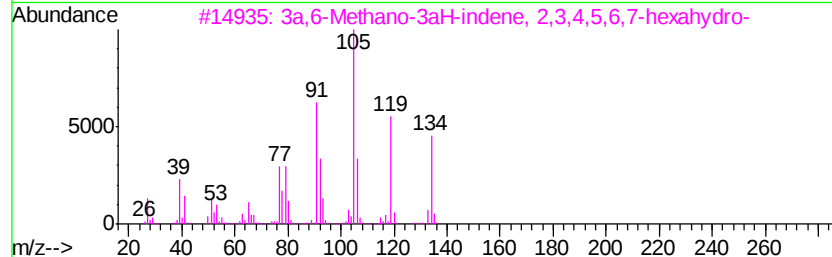
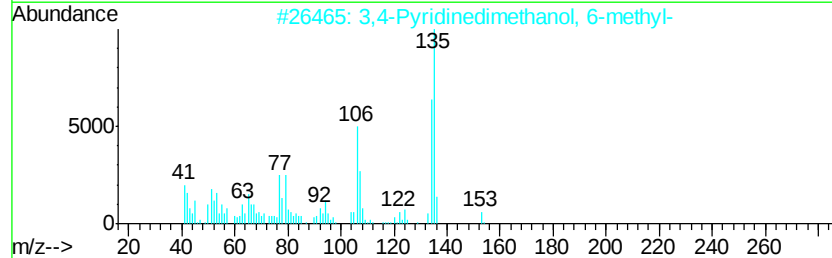
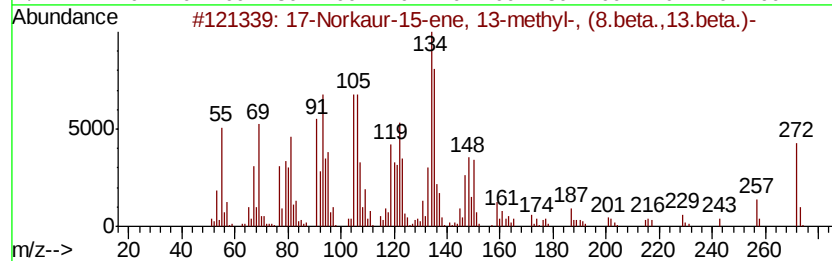
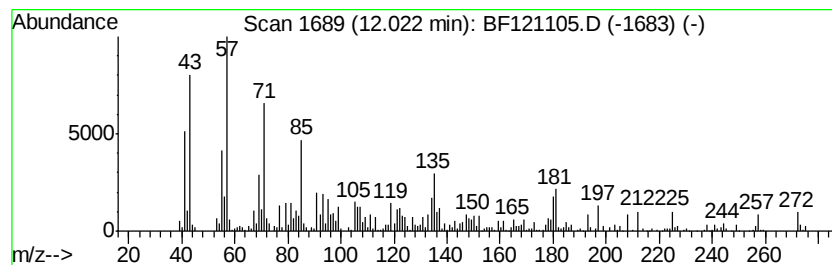
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 17-Norkaur-15-ene, 13-methy... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.02	4.32 ng	294049	Phenanthrene-d10	11.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	17-Norkaur-15-ene, 13-methyl-, (...	272 C20H32	003564-54-3	96
2	3,4-Pyridinedimethanol, 6-methyl-	153 C8H11N02	004664-11-3	55
3	3a,6-Methano-3aH-indene, 2,3,4,5...	134 C10H14	098640-10-9	35
4	Ethyltetramethylcyclopentadiene	150 C11H18	057693-77-3	35
5	2-Allylphenol	134 C9H10O	001745-81-9	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\
Data File : BF121105.D
Acq On : 29 Jul 2020 16:52
Operator : JU/CG
Sample : L3430-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

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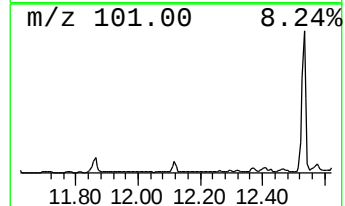
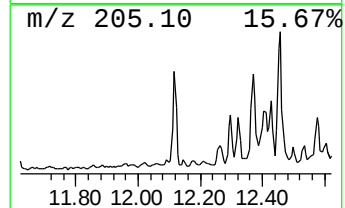
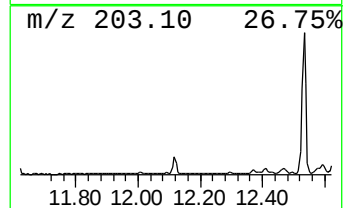
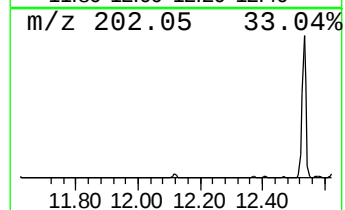
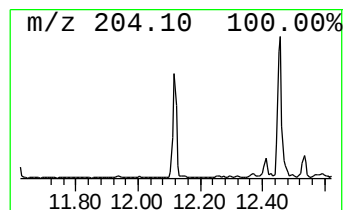
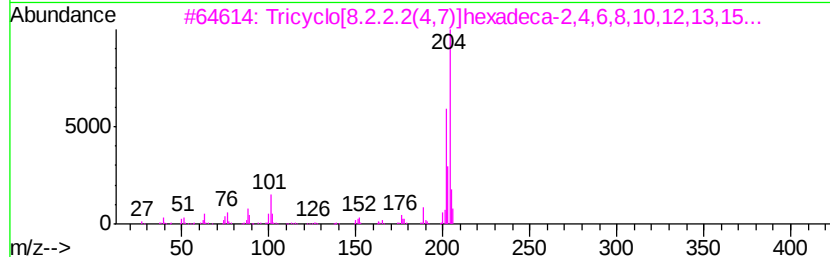
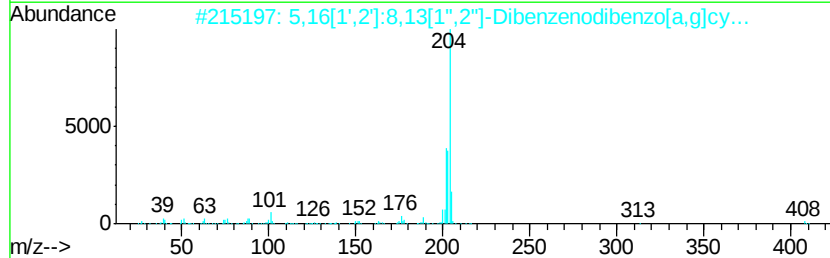
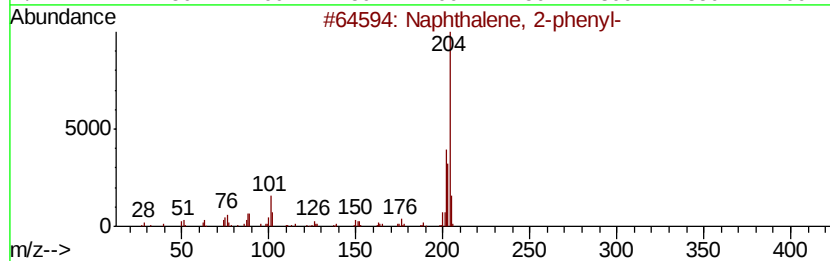
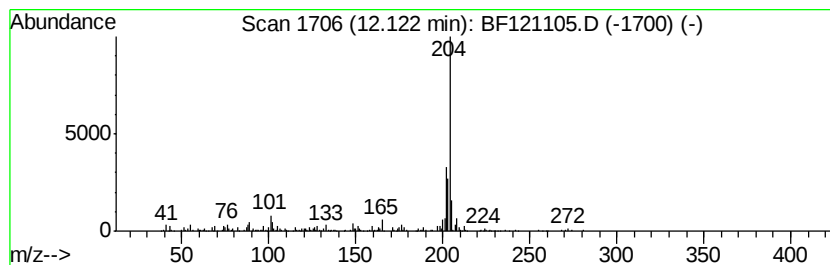
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Naphthalene, 2-phenyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.12	4.88 ng	332166	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	93
2		5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	87
3		Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	68
4		Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	58
5		9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\
Data File : BF121105.D
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Sample : L3430-02
Misc :
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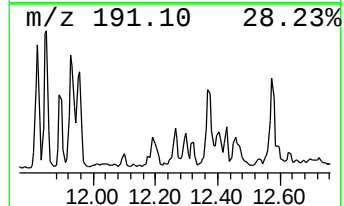
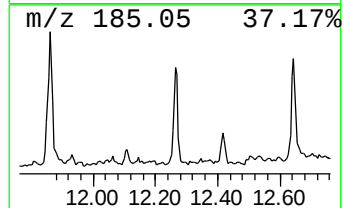
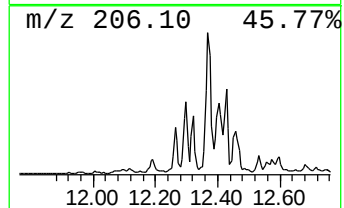
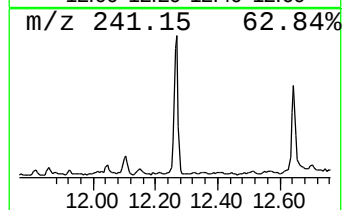
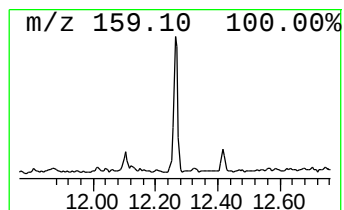
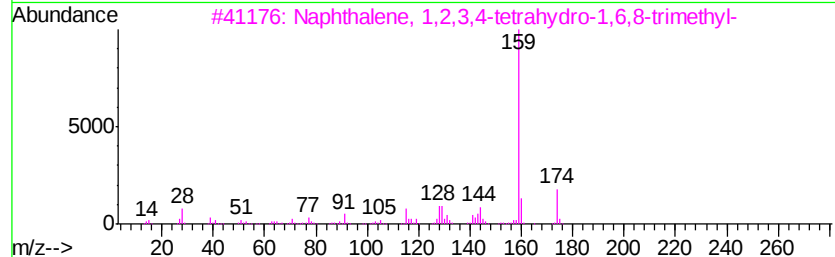
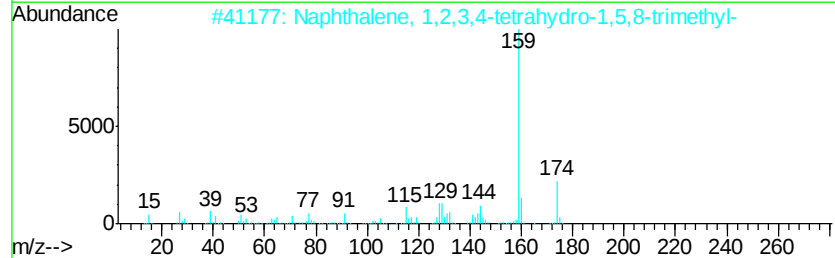
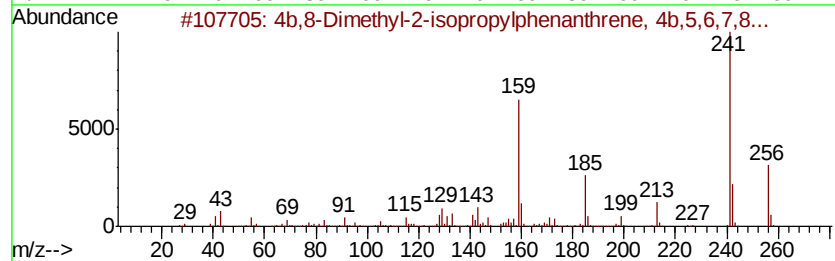
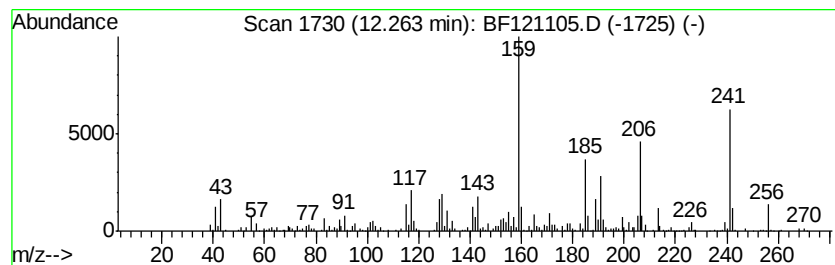
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 4b,8-Dimethyl-2-isopropylph... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.26	5.59 ng	380051	Phenanthrene-d10	11.32	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4b,8-Dimethyl-2-isopropylphenant...	256	C19H28	1000197-14-1	86
2	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	021693-51-6	30
3	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	030316-36-0	25
4	Naphthalene, 1,2,3,4-tetrahydro-...	202	C15H22	000483-77-2	25
5	Quinoline, 6-methoxy-	159	C10H9NO	005263-87-6	22



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\
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Sample : L3430-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

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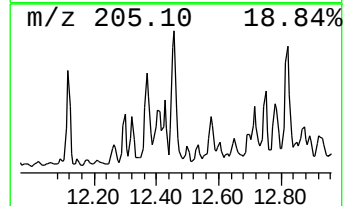
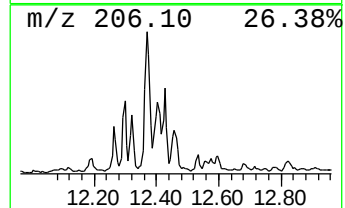
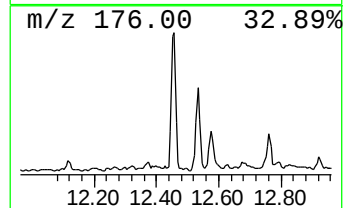
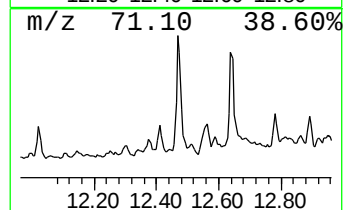
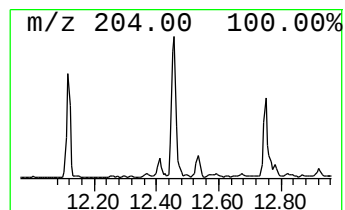
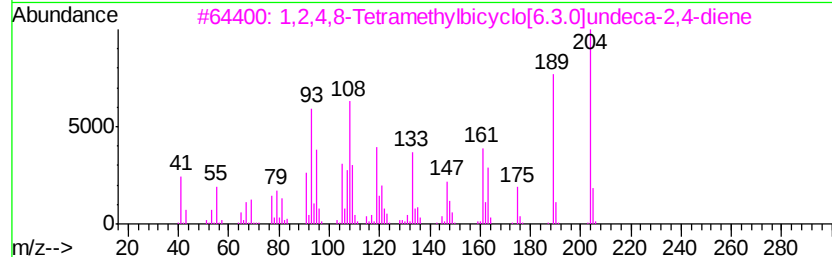
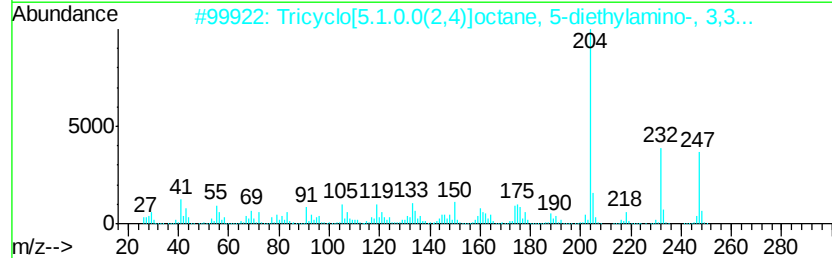
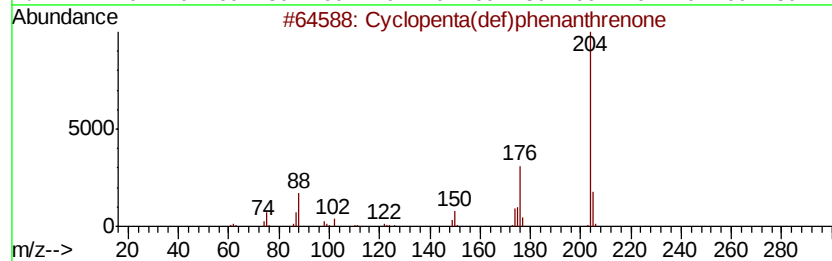
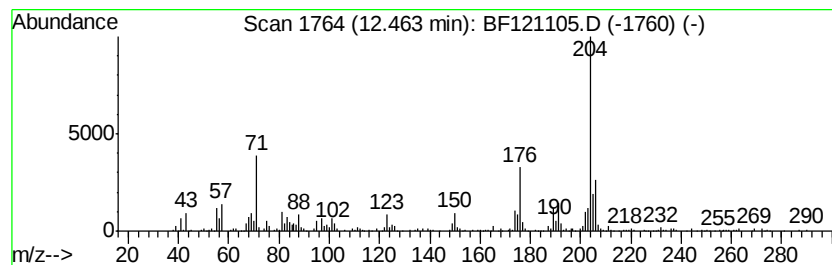
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Cyclopenta(def)phenanthrenone Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.46	4.60 ng	312723	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	94
2		Tricyclo[5.1.0.0(2,4)]octane, 5-...	247	C17H29N	1000063-59-3	59
3		1,2,4,8-Tetramethylbicyclo[6.3.0]undecane	204	C15H24	137235-51-9	53
4		[1,1'-Biphenyl]-2,2'-dicarbonitrile	204	C14H8N2	004341-02-0	50
5		2-Methyl-5-nitroindole-3-carboxamide	204	C10H8N2O3	003558-17-6	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\
Data File : BF121105.D
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Misc :
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Instrument :
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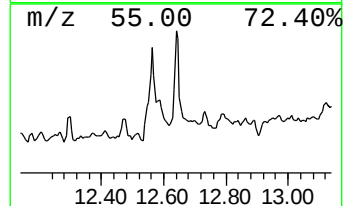
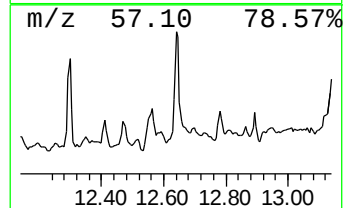
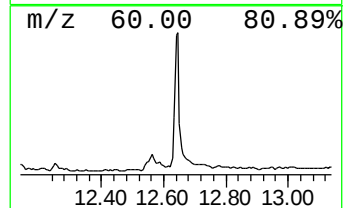
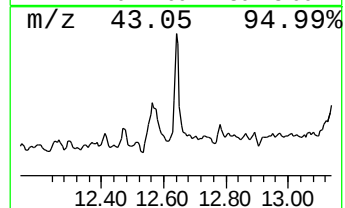
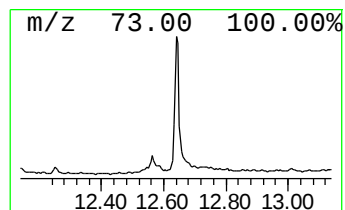
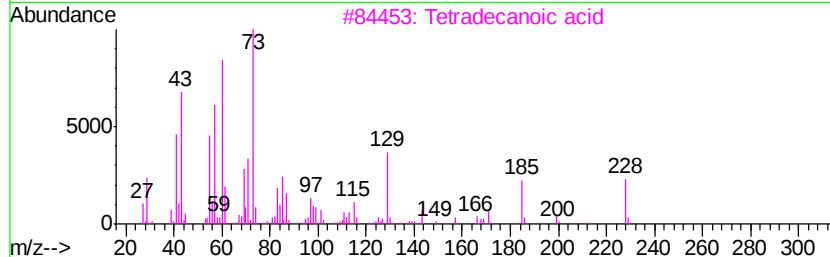
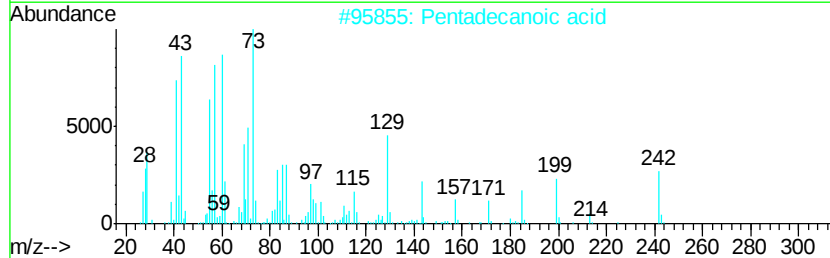
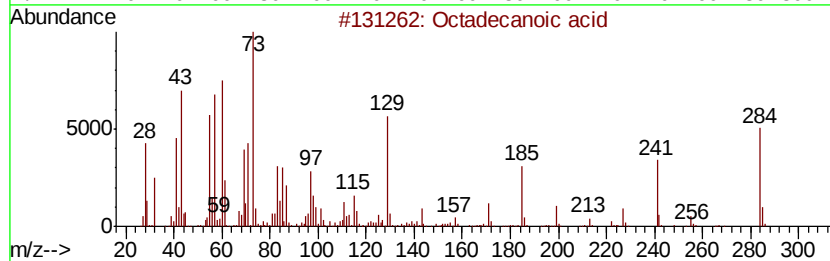
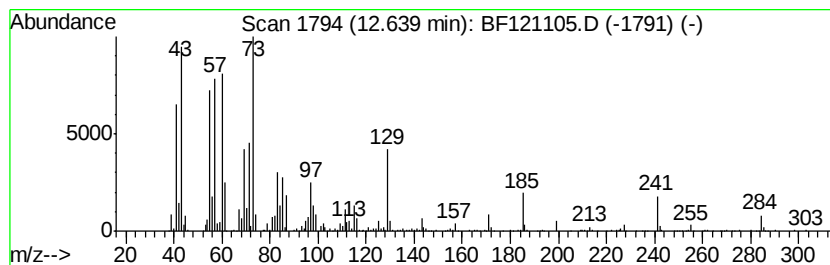
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.64	9.89 ng	673037	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	93
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	74
4		n-Decanoic acid	172	C10H20O2	000334-48-5	58
5		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\
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ALS Vial : 5 Sample Multiplier: 1

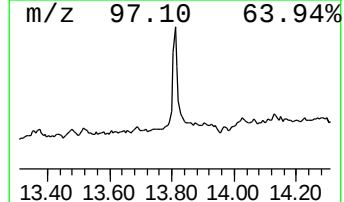
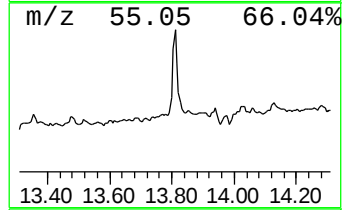
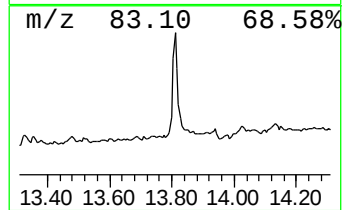
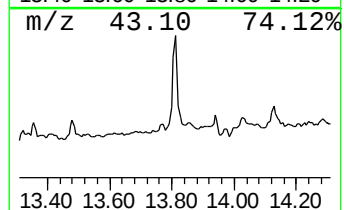
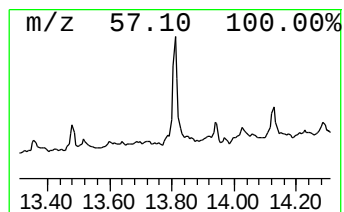
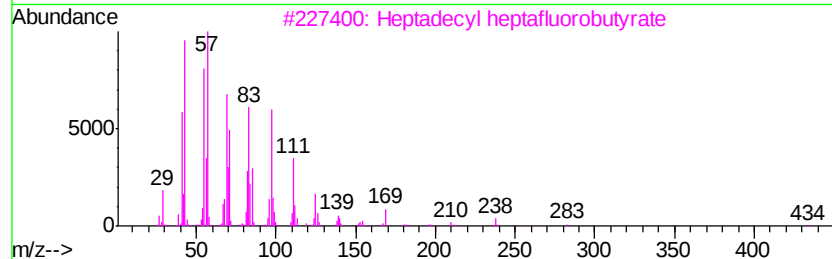
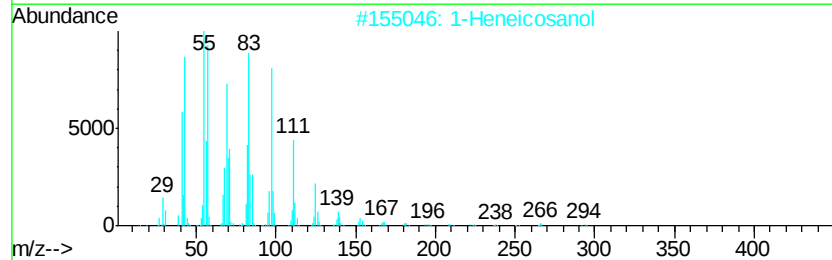
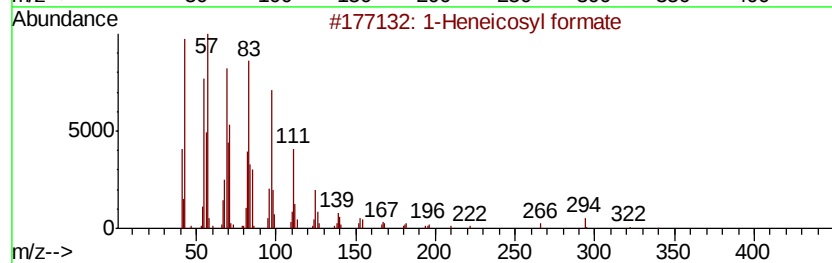
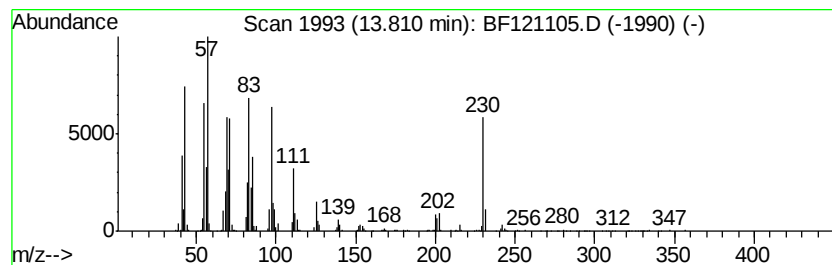
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1-Heneicosyl formate Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.81	6.52 ng	899153	Chrysene-d12		13.96	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Heneicosyl formate	340	C22H44O2	077899-03-7	95
2		1-Heneicosanol	312	C21H44O	015594-90-8	90
3		Heptadecyl heptafluorobutyrte	452	C21H35F7O2	959085-66-6	86
4		Eicosyl pentafluoropropionate	444	C23H41F5O2	1000351-80-8	76
5		Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\
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ALS Vial : 5 Sample Multiplier: 1

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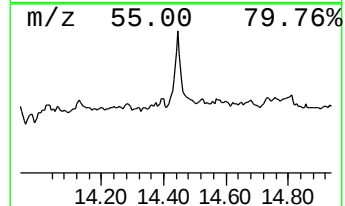
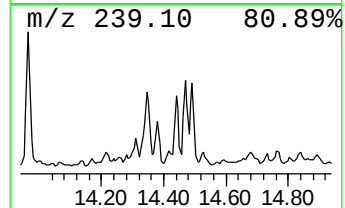
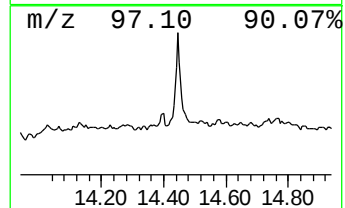
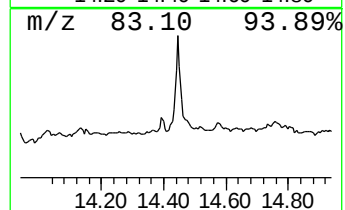
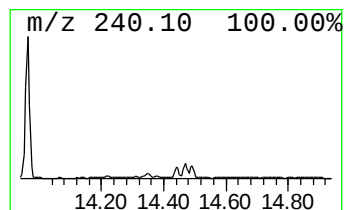
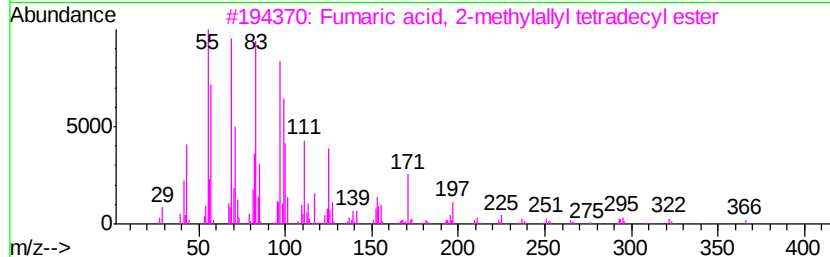
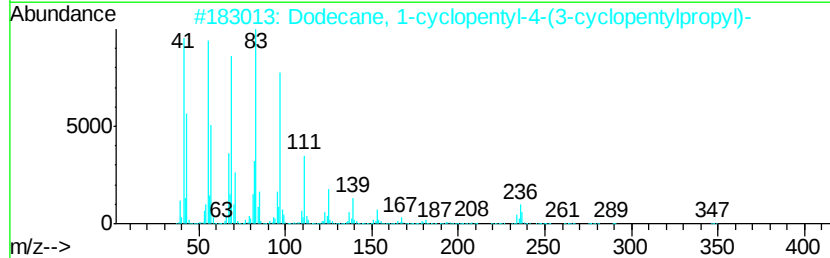
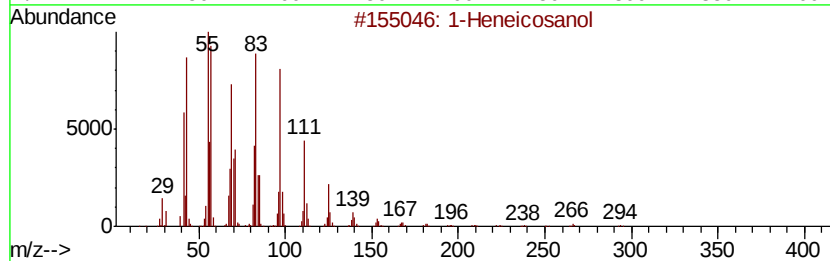
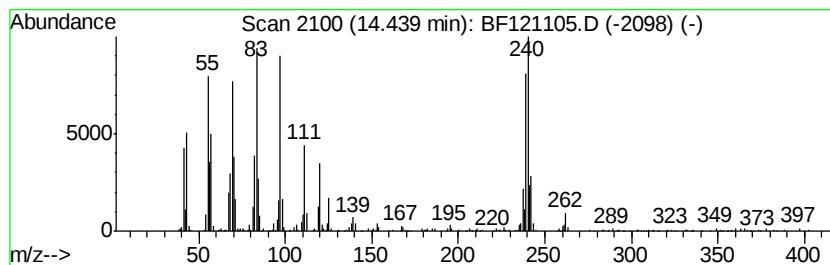
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TIC Integration Parameters: LSCINT.P

Peak Number 10 1-Heneicosanol Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.44	3.69 ng	508636	Chrysene-d12	13.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosanol	312	C21H44O	015594-90-8	74
2			Dodecane, 1-cyclopentyl-4-(3-cyclopentylpropyl)-	348	C25H48	007225-68-5	58
3			Fumaric acid, 2-methylallyl tetradecyl ester	366	C22H38O4	1000330-55-3	38
4			Cyclopentane, (2-methylbutyl)-	140	C10H20	053366-38-4	35
5			1-Octadecanol	270	C18H38O	000112-92-5	25



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Data File : BF121105.D
Acq On : 29 Jul 2020 16:52
Operator : JU/CG
Sample : L3430-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

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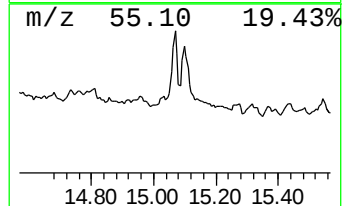
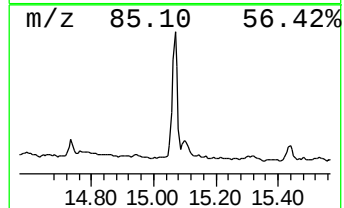
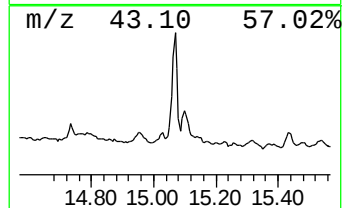
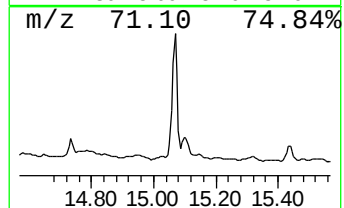
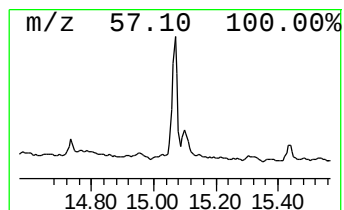
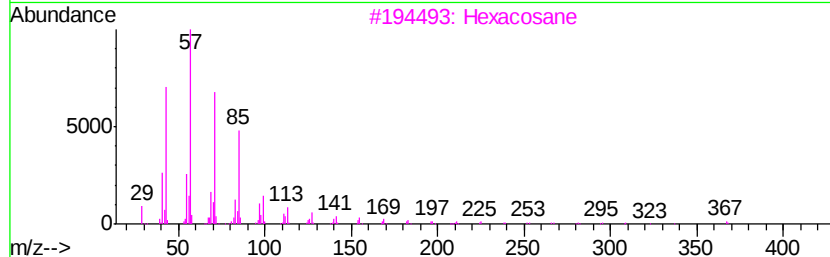
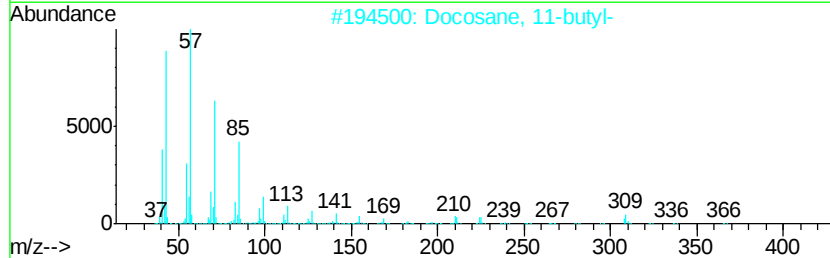
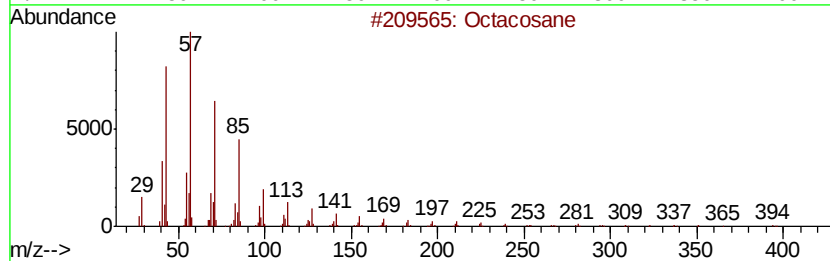
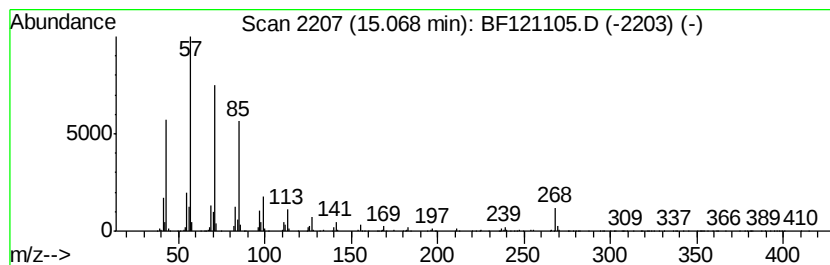
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.07	21.14 ng	1080580	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	95
2		Docosane, 11-butyl-	366	C26H54	013475-76-8	90
3		Hexacosane	366	C26H54	000630-01-3	90
4		Heneicosane	296	C21H44	000629-94-7	90
5		Heptadecane	240	C17H36	000629-78-7	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\
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ClientSampleId :
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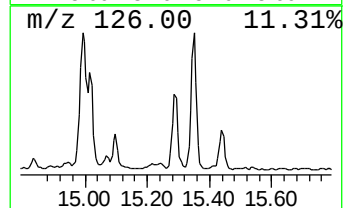
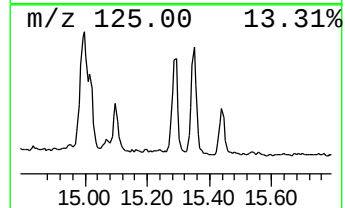
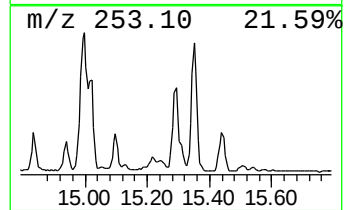
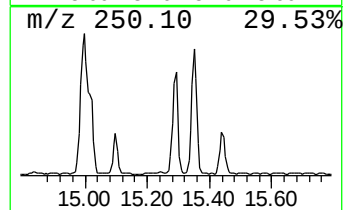
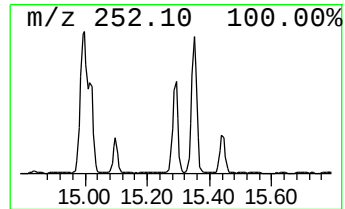
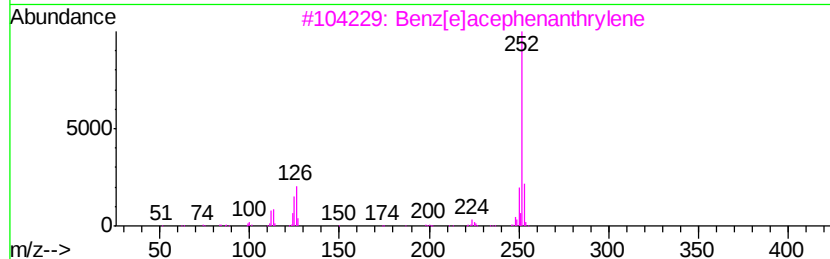
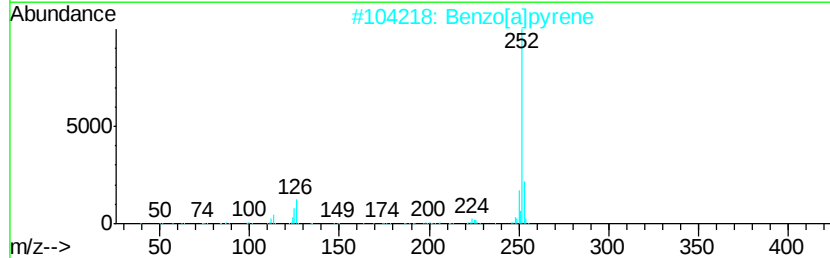
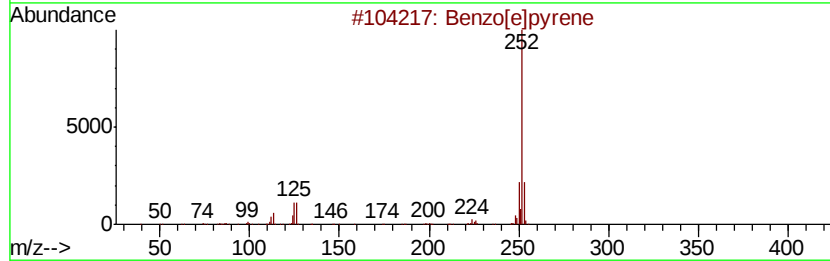
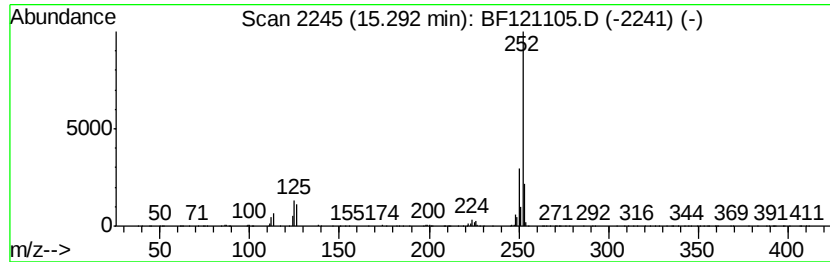
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Benzo[e]pyrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.29	18.97 ng	969593	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2		Benzo[a]pyrene	252	C20H12	000050-32-8	97
3		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	96
4		Perylene	252	C20H12	000198-55-0	93
5		Benzo[j]fluoranthene	252	C20H12	000205-82-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\
Data File : BF121105.D
Acq On : 29 Jul 2020 16:52
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Sample : L3430-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

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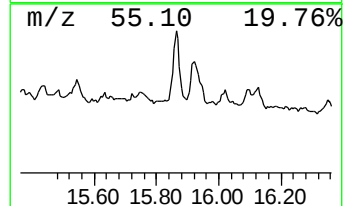
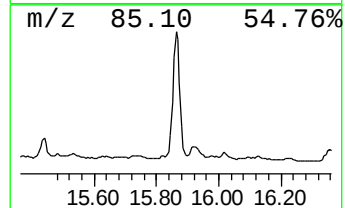
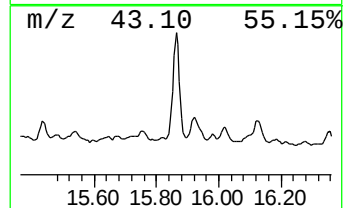
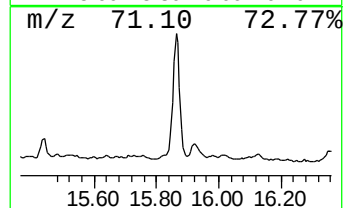
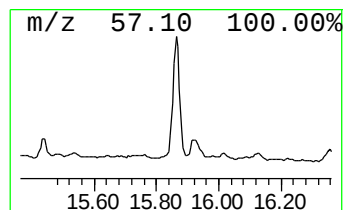
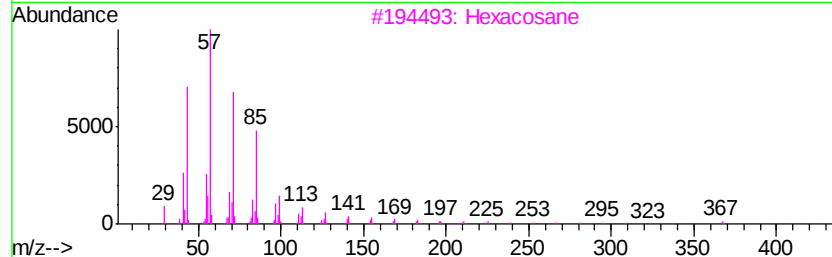
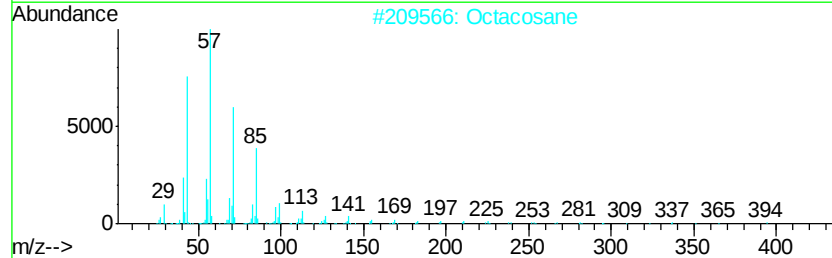
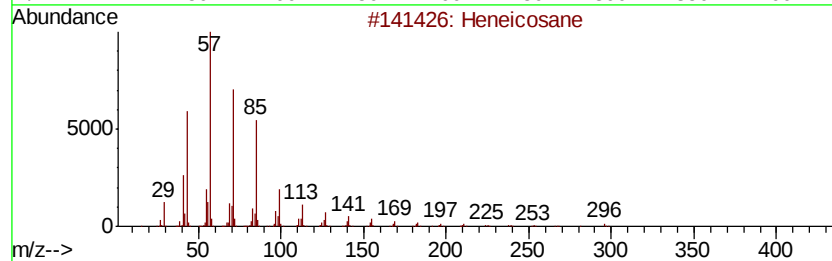
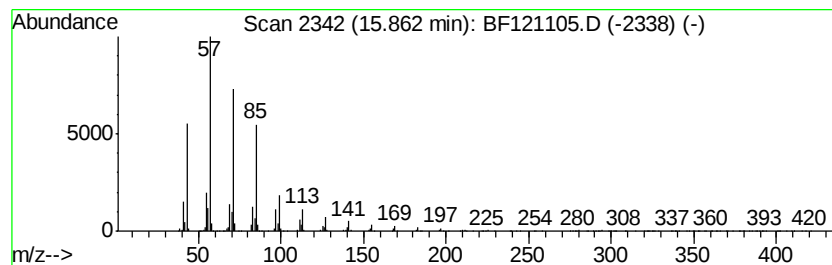
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Heneicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.86	20.86 ng	1066260	Perylene-d12	15.42

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Octacosane		394	C28H58	000630-02-4	91
3	Hexacosane		366	C26H54	000630-01-3	91
4	Tetracosane		338	C24H50	000646-31-1	91
5	Eicosane, 7-hexyl-		366	C26H54	055333-99-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\
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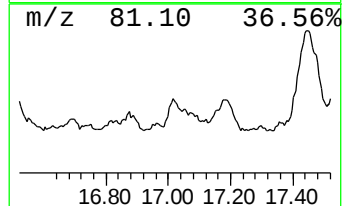
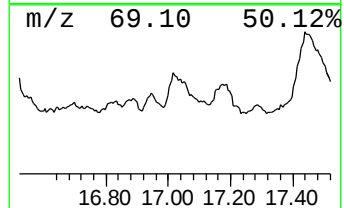
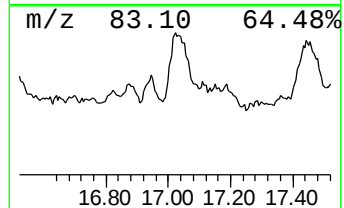
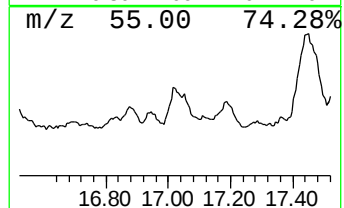
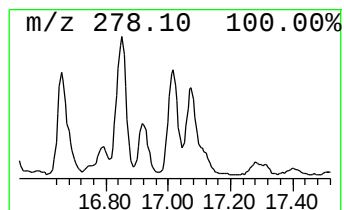
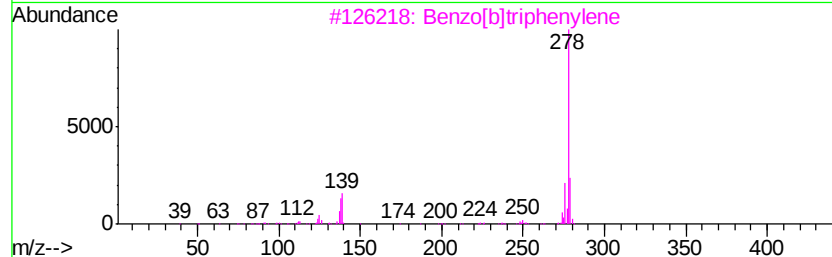
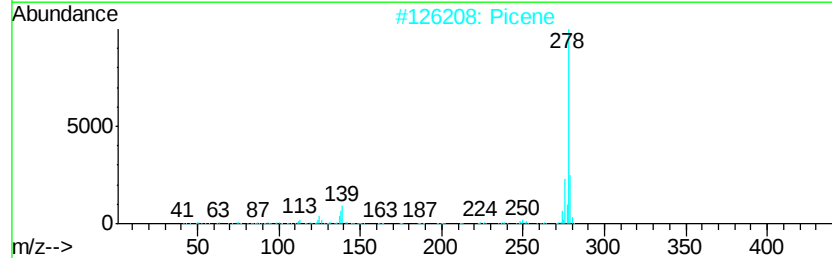
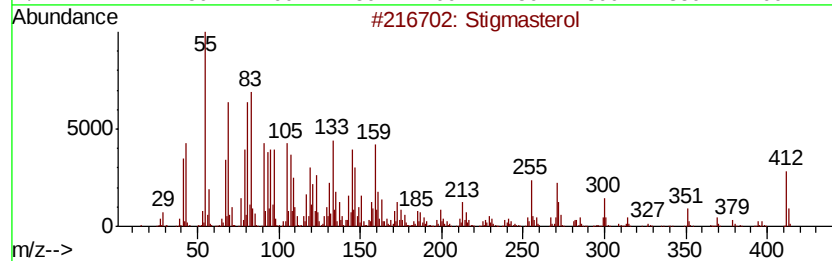
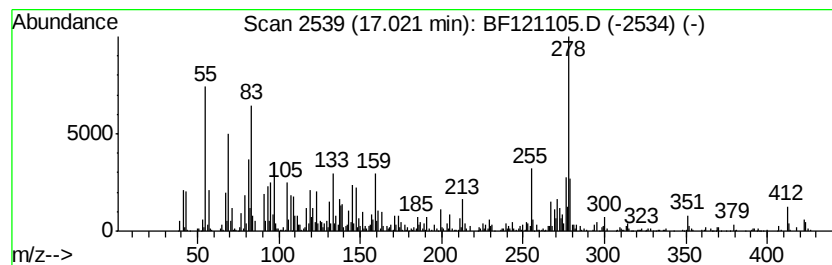
Instrument :
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Stigmasterol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.02	15.12 ng	772986	Perylene-d12	15.42
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Stigmasterol	412	C29H48O	000083-48-7 91
2	Picene	278	C22H14	000213-46-7 90
3	Benzo[b]triphenylene	278	C22H14	000215-58-7 90
4	Dibenz[a,h]anthracene	278	C22H14	000053-70-3 89
5	Quinoxaline, 6-(3-nitrobenzylidene...	278	C15H10N4O2	302800-55-1 89



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\
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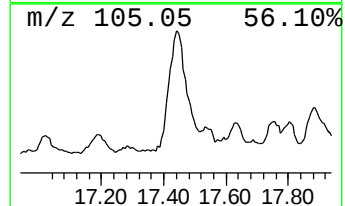
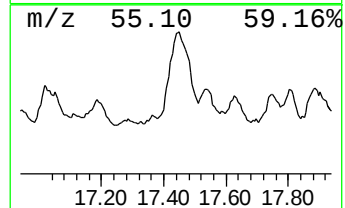
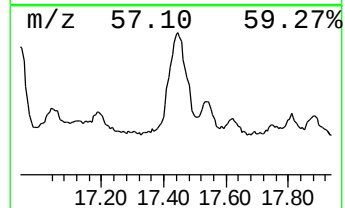
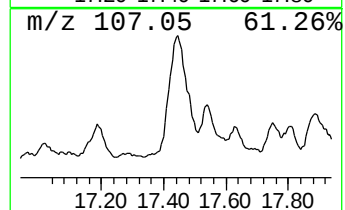
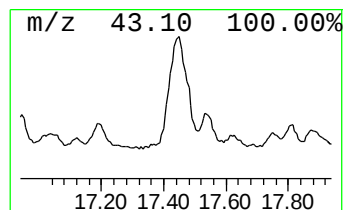
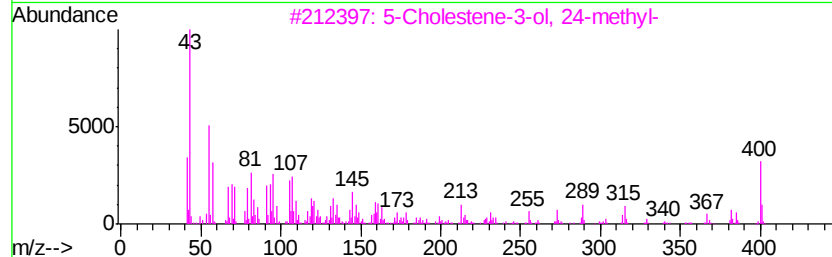
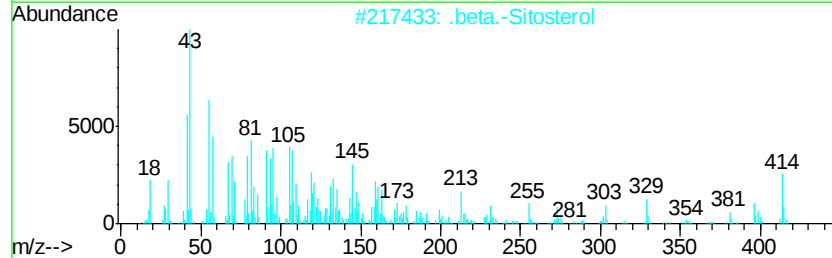
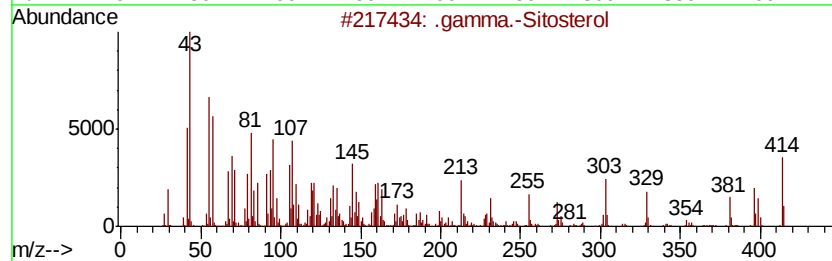
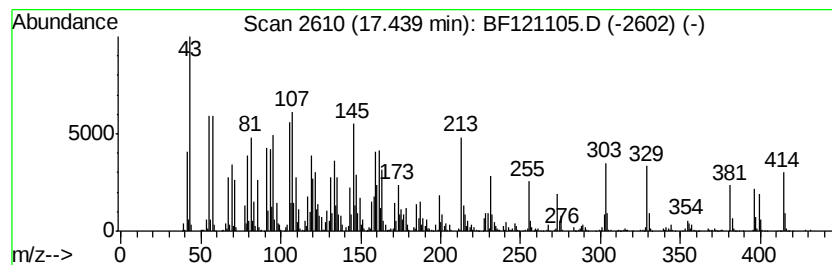
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 .gamma.-Sitosterol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.44	130.49 ng	6670000	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.gamma.-Sitosterol	414	C29H50O	000083-47-6	99
2		.beta.-Sitosterol	414	C29H50O	000083-46-5	89
3		5-Cholestene-3-ol, 24-methyl-	400	C28H48O	1000214-17-4	38
4		22,26-Oxido-4,17-cholestadien-3....	414	C27H42O3	1000252-80-4	11
5		Ergost-5-en-3-ol, (3.beta.)-	400	C28H48O	004651-51-8	11



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\
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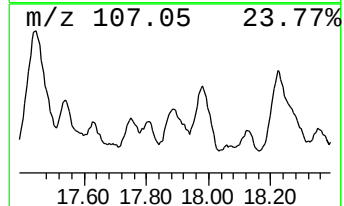
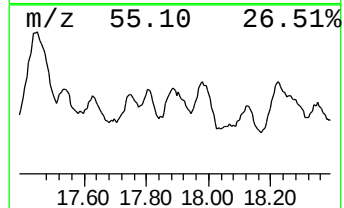
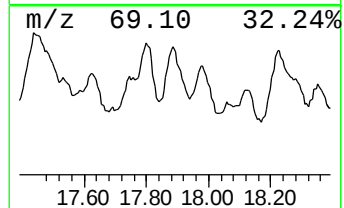
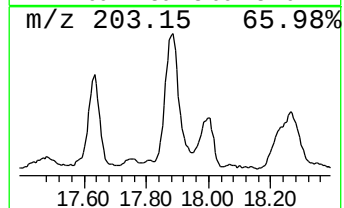
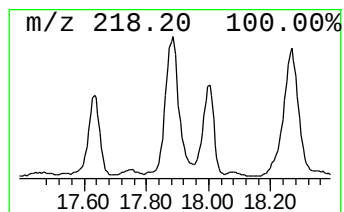
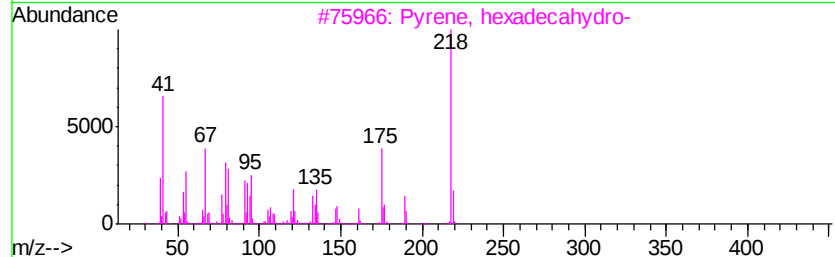
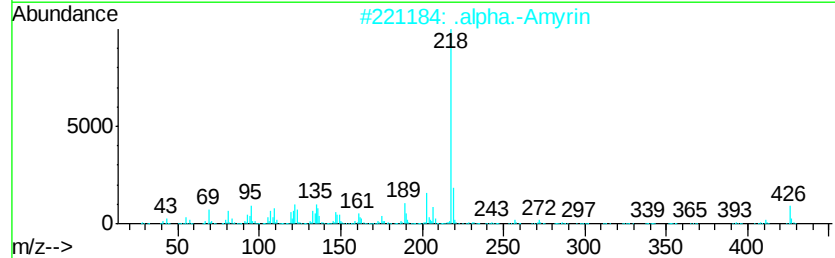
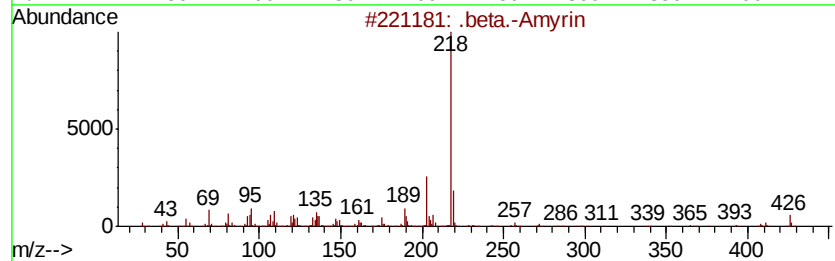
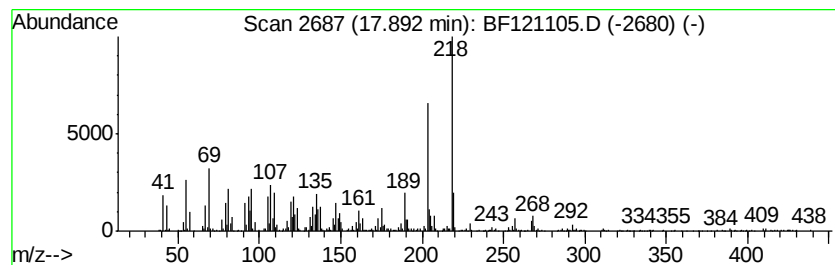
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 .beta.-Amyrin Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.89	38.18 ng	1951840	Perylene-d12	15.42

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		.beta.-Amyrin	426	C30H500	000559-70-6	92
2		.alpha.-Amyrin	426	C30H500	000638-95-9	78
3		Pvrene, hexadecahydro-	218	C16H26	002435-85-0	55
4		4,4,6a,6b,8a,11,11,14b-Octamethy...	424	C30H480	1000194-62-4	52
5		3-Methoxy-6,7,8,9-tetrahydro-dib...	218	C13H1403	1000186-57-9	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\
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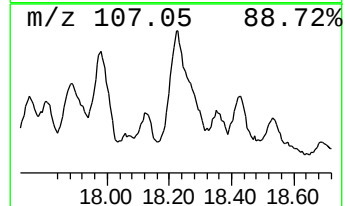
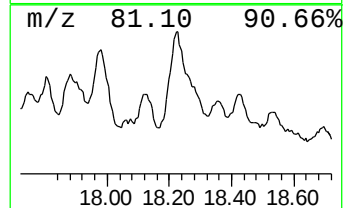
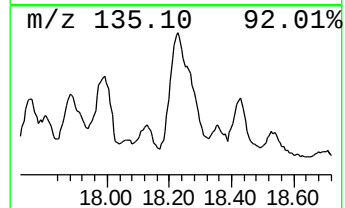
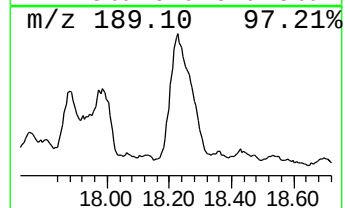
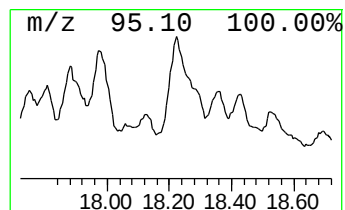
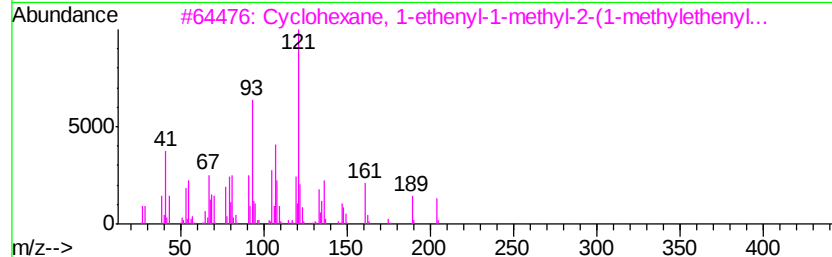
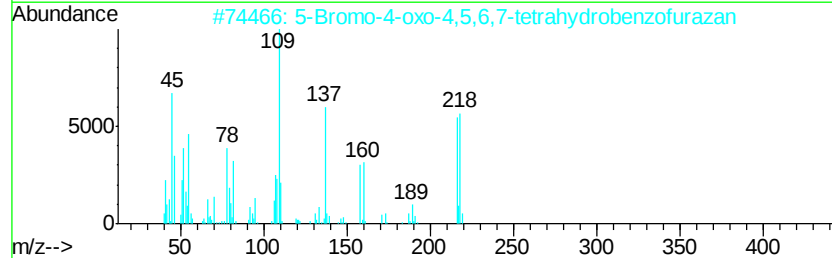
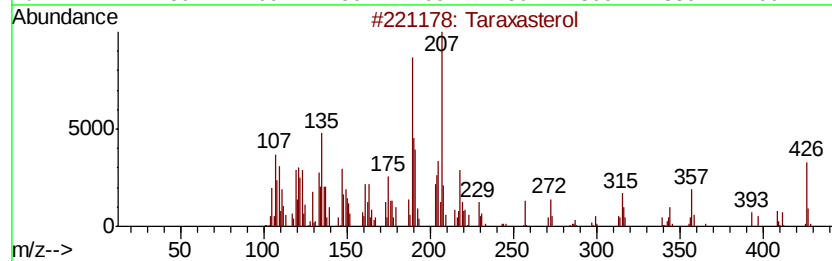
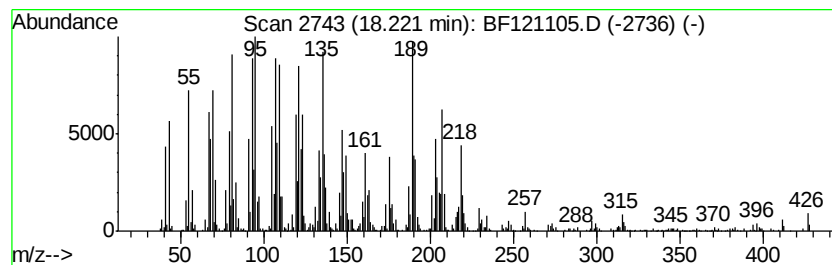
Instrument :
BNA_F
ClientSampleId :
TS2

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 19 Taraxasterol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.22	88.53 ng	4525020	Perylene-d12	15.42
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Taraxasterol	426 C30H50O	001059-14-9 72
2		5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216 C6H5BrN2O2	300574-36-1 59
3		Cyclohexane, 1-ethenyl-1-methyl-...	204 C15H24	003242-08-8 46
4		4,6,6-Trimethyl-2-(3-methylbuta-...	218 C15H22O	1000190-22-2 45
5		7-Oxabicyclo[4.1.0]heptane, 2,2,...	218 C15H22O	070038-20-9 44



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\
Data File : BF121105.D
Acq On : 29 Jul 2020 16:52
Operator : JU/CG
Sample : L3430-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TS2

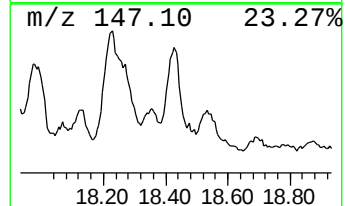
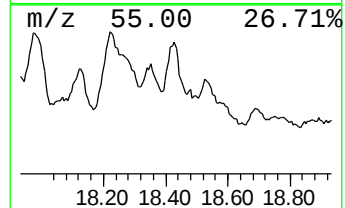
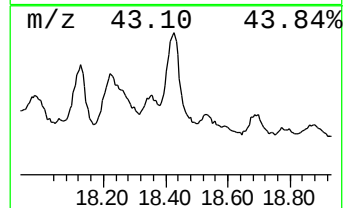
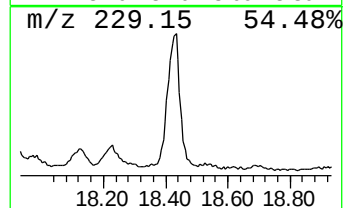
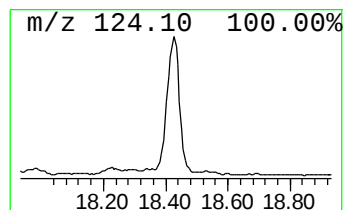
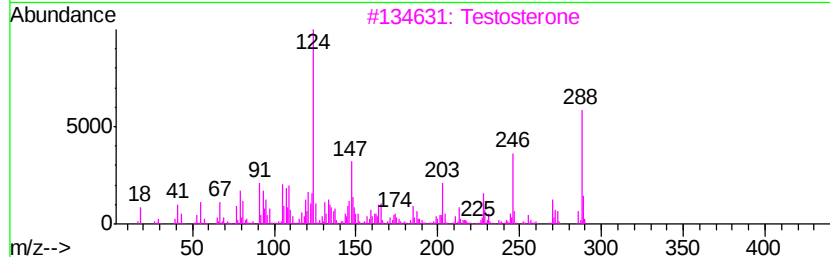
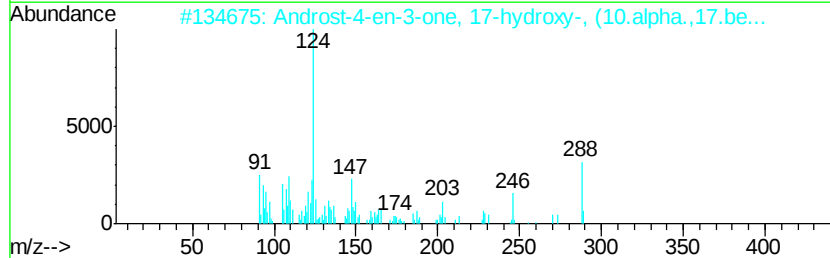
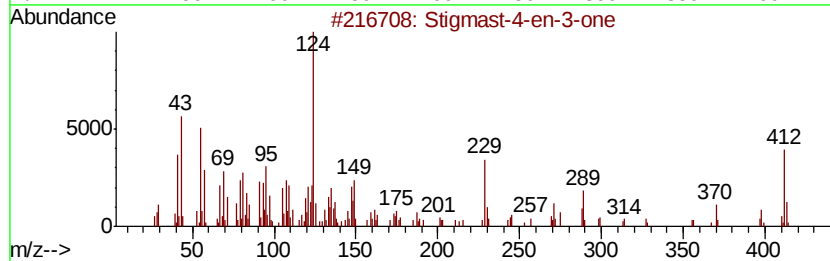
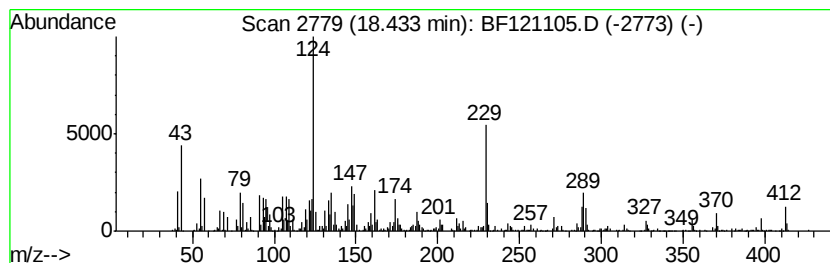
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Stigmast-4-en-3-one Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.43	41.63 ng	2127900	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Stigmast-4-en-3-one	412	C29H48O	001058-61-3	60
2		Androst-4-en-3-one, 17-hydroxy-, ...	288	C19H28O2	000604-39-7	50
3		Testosterone	288	C19H28O2	000058-22-0	43
4		Pregn-4-ene-3,20-dione, (9.beta....	314	C21H30O2	002755-10-4	43
5		Testosterone cypionate	412	C27H40O3	000058-20-8	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF072920\
 Data File : BF121105.D
 Acq On : 29 Jul 2020 16:52
 Operator : JU/CG
 Sample : L3430-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS2

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
Anthracene, 2-met...	11.82	5.5 ng		373700	4	11.32	1360940	20.0
n-Hexadecanoic acid	11.86	24.9 ng		1692610	4	11.32	1360940	20.0
4H-Cyclopenta[def...	11.93	3.7 ng		253561	4	11.32	1360940	20.0
17-Norkaur-15-ene...	12.02	4.3 ng		294049	4	11.32	1360940	20.0
Naphthalene, 2-ph...	12.12	4.9 ng		332166	4	11.32	1360940	20.0
4b,8-Dimethyl-2-i...	12.26	5.6 ng		380051	4	11.32	1360940	20.0
Cyclopenta(def)ph...	12.46	4.6 ng		312723	4	11.32	1360940	20.0
Octadecanoic acid	12.64	9.9 ng		673037	4	11.32	1360940	20.0
1-Heneicosyl formate	13.81	6.5 ng		899153	5	13.96	2756520	20.0
1-Heneicosanol	14.44	3.7 ng		508636	5	13.96	2756520	20.0
Octacosane	15.07	21.1 ng		1080580	6	15.42	1022310	20.0
Benzo[e]pyrene	15.29	19.0 ng		969593	6	15.42	1022310	20.0
Heneicosane	15.86	20.9 ng		1066260	6	15.42	1022310	20.0
Stigmasterol	17.02	15.1 ng		772986	6	15.42	1022310	20.0
.gamma.-Sitosterol	17.44	130.5 ng		6670000	6	15.42	1022310	20.0
.beta.-Amyrin	17.89	38.2 ng		1951840	6	15.42	1022310	20.0
4,4,6a,6b,8a,11,1...	17.98	47.4 ng		2422360	6	15.42	1022310	20.0
Taraxasterol	18.22	88.5 ng		4525020	6	15.42	1022310	20.0
Stigmast-4-en-3-one	18.43	41.6 ng		2127900	6	15.42	1022310	20.0