

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071919\
 Data File : BF115670.D
 Acq On : 19 Jul 2019 12:07
 Operator : HP/JU
 Sample : K3618-05 2X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HD-01-071719

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-BF071219.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	2.169	9	14	20	rVB	322591	421271	19.82%	1.388%
2	5.498	576	580	593	rBV	1792100	1663834	78.27%	5.483%
3	6.504	747	751	758	rBV	1800424	1788140	84.12%	5.892%
4	6.557	758	760	764	rVB	36250	35955	1.69%	0.118%
5	6.651	772	776	779	rBV	2136197	1869292	87.94%	6.160%
6	6.881	811	815	818	rBV	1543870	1255113	59.04%	4.136%
7	7.033	837	841	845	rBV	1559010	1461749	68.76%	4.817%
8	7.263	877	880	890	rBV	17627	31327	1.47%	0.103%
9	7.433	905	909	921	rBV	1227315	1164517	54.78%	3.837%
10	8.163	1029	1033	1036	rBV	1985315	1616912	76.06%	5.328%
11	8.222	1040	1043	1057	rVB4	20369	43630	2.05%	0.144%
12	8.975	1167	1171	1177	rVB	25925	22849	1.07%	0.075%
13	9.239	1211	1216	1219	rBV	2276409	2125753	100.00%	7.005%
14	9.574	1269	1273	1275	rBV	45055	41692	1.96%	0.137%
15	9.604	1275	1278	1279	rVV	74882	55841	2.63%	0.184%
16	9.627	1279	1282	1290	rVB	424043	350604	16.49%	1.155%
17	9.692	1290	1293	1298	rBV	23403	23870	1.12%	0.079%
18	9.763	1302	1305	1306	rBV	80909	65257	3.07%	0.215%
19	9.916	1325	1331	1335	rBV	1938217	1819536	85.59%	5.996%
20	9.951	1335	1337	1340	rVB	43243	35026	1.65%	0.115%
21	10.121	1362	1366	1369	rBV2	44278	46408	2.18%	0.153%
22	10.157	1369	1372	1377	rVB	28743	25529	1.20%	0.084%
23	10.233	1381	1385	1388	rBV2	25485	25116	1.18%	0.083%
24	10.327	1397	1401	1403	rBV2	20195	29393	1.38%	0.097%
25	10.433	1416	1419	1422	rVV	96337	72303	3.40%	0.238%
26	10.463	1422	1424	1428	rVB	53824	46954	2.21%	0.155%
27	10.704	1461	1465	1476	rVB	1394026	1254953	59.04%	4.135%
28	10.827	1482	1486	1493	rVB3	41908	51326	2.41%	0.169%
29	11.016	1513	1518	1520	rBV3	27428	33865	1.59%	0.112%
30	11.039	1520	1522	1525	rBV2	25163	21854	1.03%	0.072%
31	11.204	1547	1550	1554	rVB	48769	45208	2.13%	0.149%
32	11.298	1564	1566	1571	rVB	69170	63539	2.99%	0.209%
33	11.351	1572	1575	1580	rBV6	13593	22666	1.07%	0.075%
34	11.404	1580	1584	1586	rVV	2037049	1706592	80.28%	5.624%

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 Integrator: RTE
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Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF071219.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.427	1586	1588	1591	rVV	683153	530870	24.97%	1.749%
36	11.457	1591	1593	1595	rVV	104191	90151	4.24%	0.297%
37	11.474	1595	1596	1599	rVB	71921	51232	2.41%	0.169%
38	11.510	1599	1602	1606	rBV4	34079	43911	2.07%	0.145%
39	11.627	1618	1622	1625	rBV	38979	49708	2.34%	0.164%
40	11.698	1631	1634	1636	rVB	42281	34339	1.62%	0.113%
41	11.733	1637	1640	1645	rVB	83972	71781	3.38%	0.237%
42	11.815	1651	1654	1657	rVB	41543	44102	2.07%	0.145%
43	11.857	1657	1661	1664	rBV2	32598	38016	1.79%	0.125%
44	11.904	1664	1669	1671	rVV	202789	207917	9.78%	0.685%
45	11.927	1671	1673	1680	rVV2	518605	530807	24.97%	1.749%
46	12.015	1684	1688	1690	rVV2	209305	234133	11.01%	0.772%
47	12.039	1690	1692	1697	rVB	148496	120098	5.65%	0.396%
48	12.104	1697	1703	1706	rBV3	46163	63317	2.98%	0.209%
49	12.139	1706	1709	1711	rVB	37814	33523	1.58%	0.110%
50	12.198	1716	1719	1721	rVV	110042	102979	4.84%	0.339%
51	12.221	1721	1723	1730	rVB2	107824	122229	5.75%	0.403%
52	12.327	1739	1741	1743	rBV	35038	35153	1.65%	0.116%
53	12.351	1743	1745	1747	rVV	58849	50864	2.39%	0.168%
54	12.380	1747	1750	1752	rVV	82051	67924	3.20%	0.224%
55	12.404	1752	1754	1756	rVB2	52388	42514	2.00%	0.140%
56	12.457	1759	1763	1765	rBV	161479	155668	7.32%	0.513%
57	12.492	1765	1769	1771	rVV2	86889	99781	4.69%	0.329%
58	12.539	1774	1777	1782	rBV3	84585	103087	4.85%	0.340%
59	12.615	1786	1790	1795	rBV	661426	540481	25.43%	1.781%
60	12.657	1795	1797	1799	rBV	34889	32390	1.52%	0.107%
61	12.680	1799	1801	1803	rVB	36207	30713	1.44%	0.101%
62	12.710	1803	1806	1809	rBV2	128700	129518	6.09%	0.427%
63	12.768	1814	1816	1819	rVV2	55612	67698	3.18%	0.223%
64	12.845	1823	1829	1832	rVV	673751	753590	35.45%	2.483%
65	12.904	1836	1839	1842	rVB2	67886	80130	3.77%	0.264%
66	12.986	1849	1853	1856	rBV	1753351	1530799	72.01%	5.044%
67	13.062	1862	1866	1870	rBV2	86365	129768	6.10%	0.428%
68	13.121	1873	1876	1878	rBV2	36547	39936	1.88%	0.132%
69	13.151	1878	1881	1882	rVV2	65910	69525	3.27%	0.229%
70	13.168	1882	1884	1887	rVB	107989	108420	5.10%	0.357%
71	13.221	1890	1893	1894	rBV	53640	47885	2.25%	0.158%

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72	13.239	1894	1896	1899	rVV	48186	42965	2.02%	0.142%
73	13.274	1899	1902	1907	rVV2	103960	107648	5.06%	0.355%
74	13.368	1916	1918	1921	rVB	92293	67473	3.17%	0.222%
75	13.398	1921	1923	1927	rVB	71950	64755	3.05%	0.213%
76	13.680	1968	1971	1976	rVB	87506	110617	5.20%	0.365%
77	13.886	2004	2006	2011	rBV2	78414	69872	3.29%	0.230%
78	14.045	2028	2033	2035	rBV2	1145270	1315054	61.86%	4.334%
79	14.068	2035	2037	2040	rVB	216104	175240	8.24%	0.577%
80	14.204	2058	2060	2067	rVB4	88525	116975	5.50%	0.385%
81	14.509	2110	2112	2117	rVB2	104757	111963	5.27%	0.369%
82	14.898	2175	2178	2182	rVB	305448	275471	12.96%	0.908%
83	15.086	2207	2210	2217	rVV	176062	318351	14.98%	1.049%
84	15.162	2220	2223	2226	rVB	143669	154401	7.26%	0.509%
85	15.386	2259	2261	2265	rVB	95039	104751	4.93%	0.345%
86	15.451	2270	2272	2276	rVB	124129	132846	6.25%	0.438%
87	15.521	2279	2284	2287	rBV	943972	1152382	54.21%	3.797%
88	15.986	2360	2363	2370	rVB	140340	202560	9.53%	0.667%

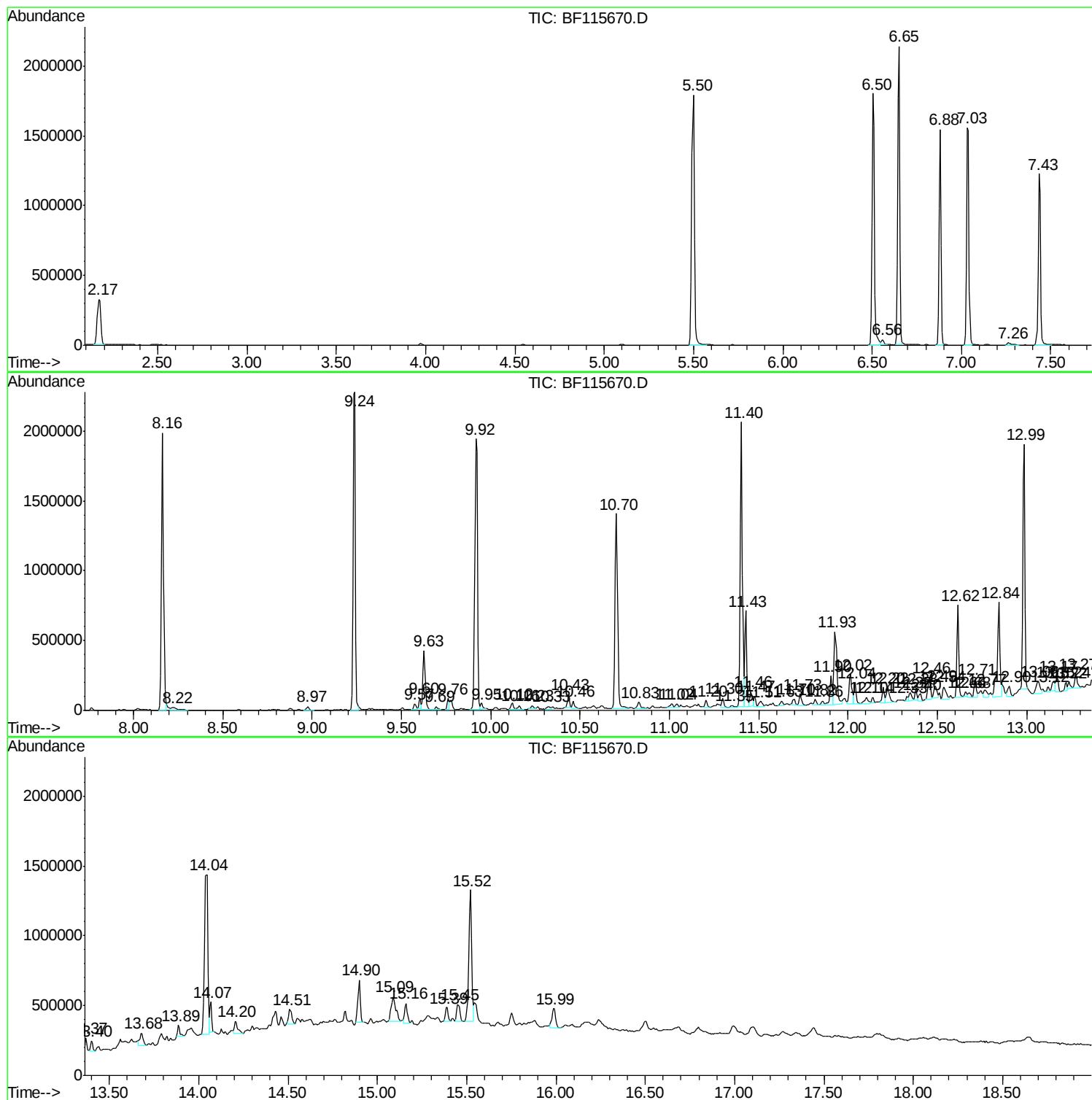
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Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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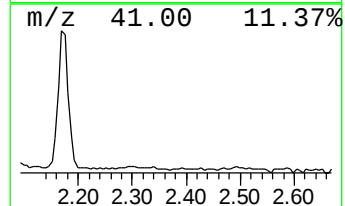
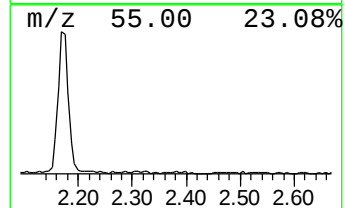
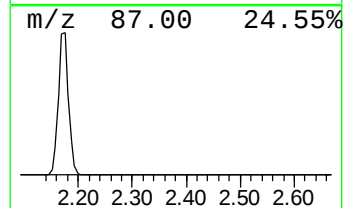
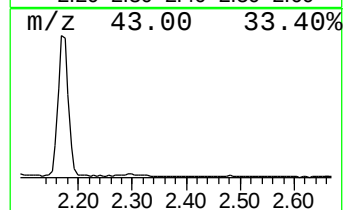
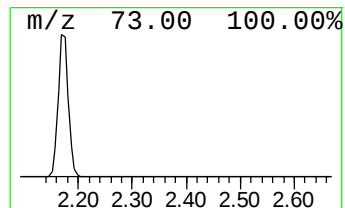
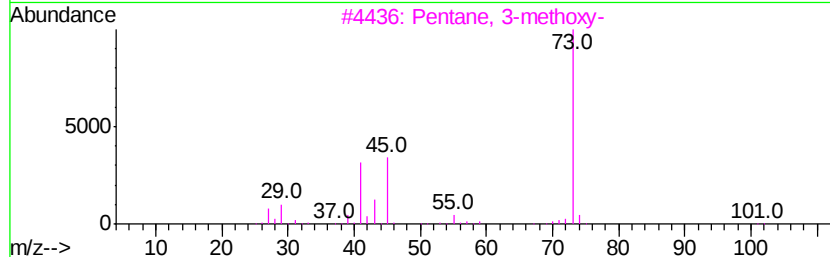
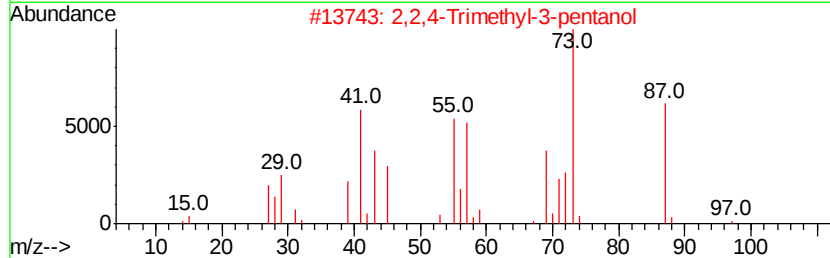
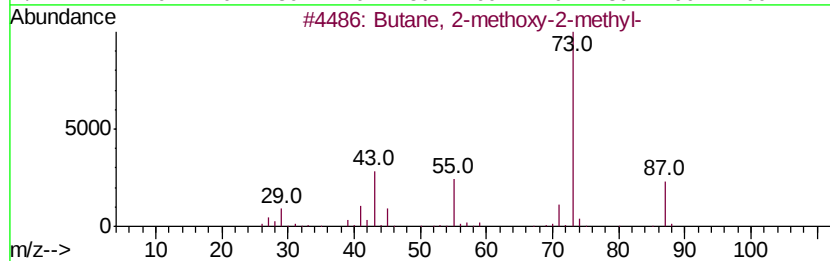
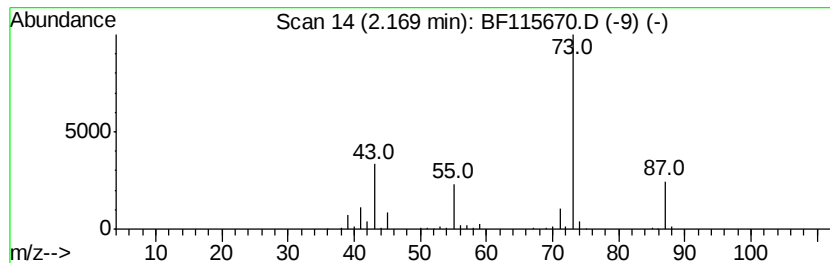
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.17	6.71 ng	421271	1,4-Dichlorobenzene-d4	6.88

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	10
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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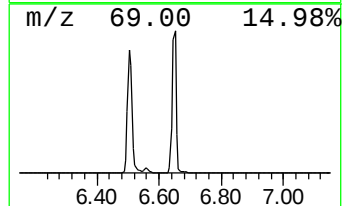
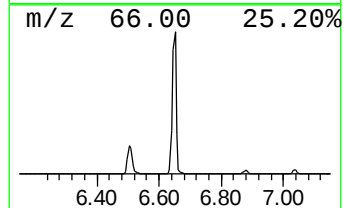
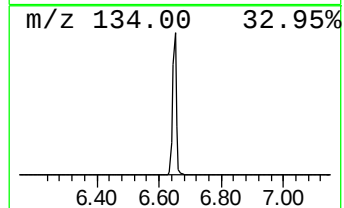
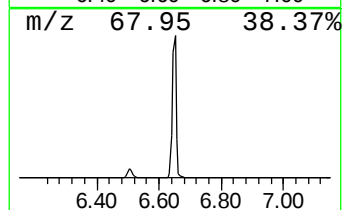
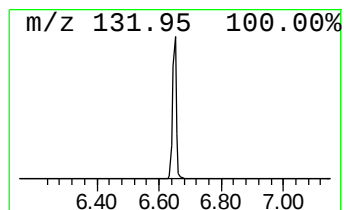
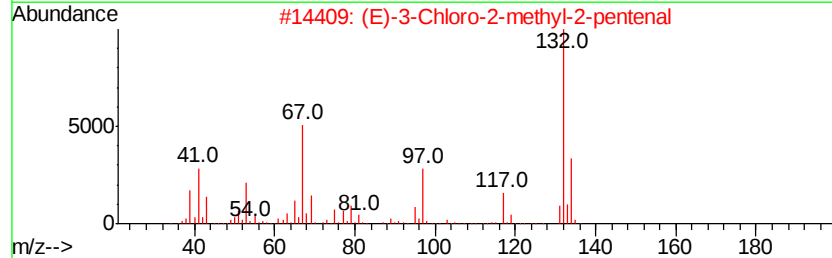
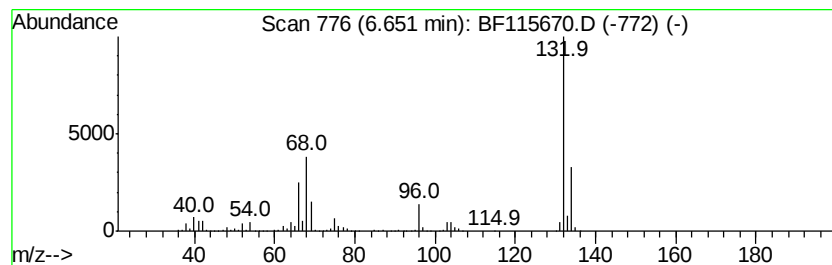
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	29.79 ng	1869290	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	12
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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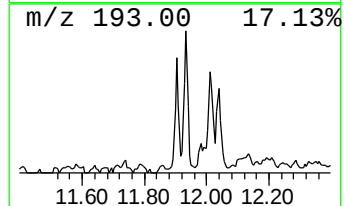
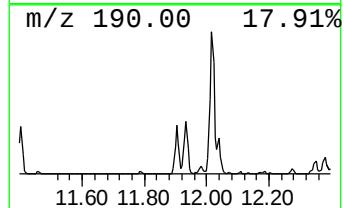
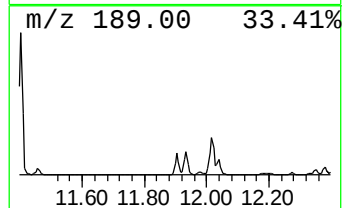
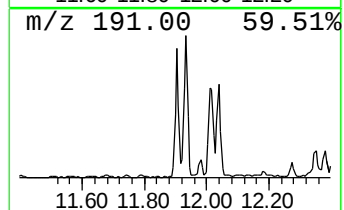
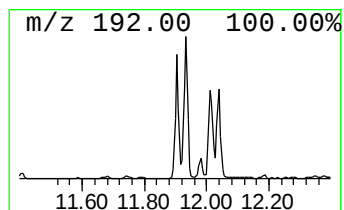
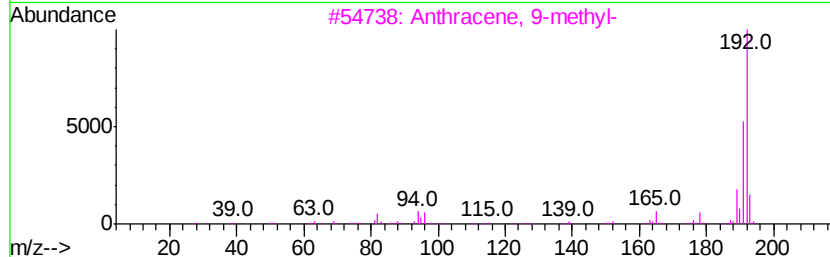
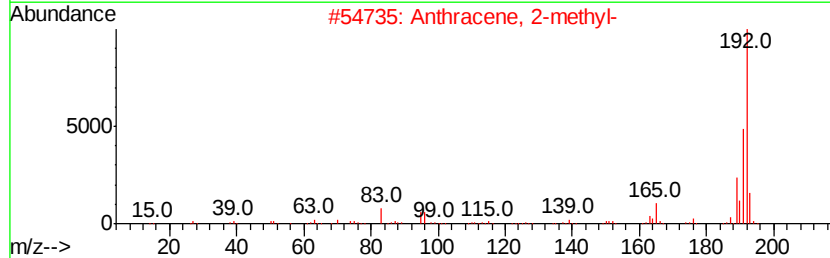
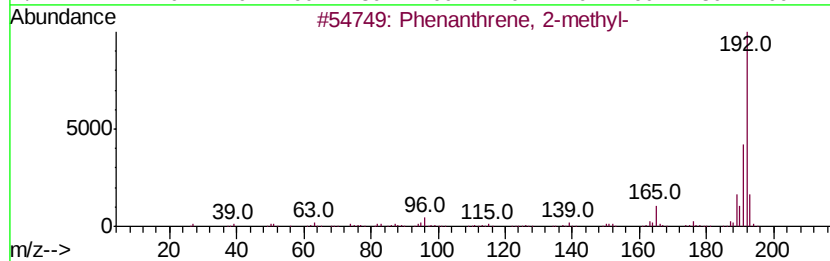
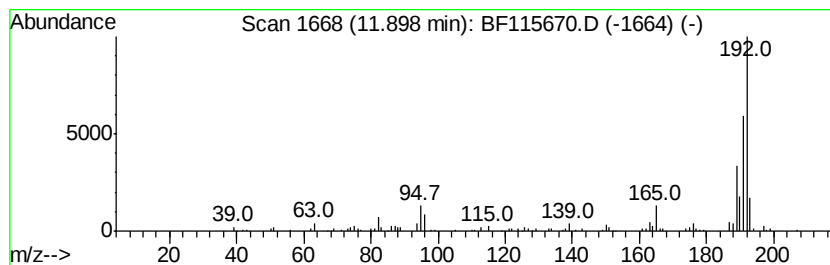
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Phenanthrene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.90	2.44 ng	207917	Phenanthrene-d10	11.40	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
3	Anthracene, 9-methyl-	192	C15H12	000779-02-2	93
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	91



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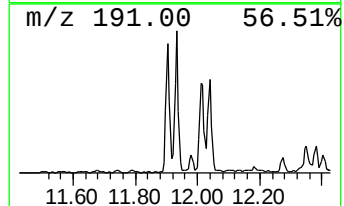
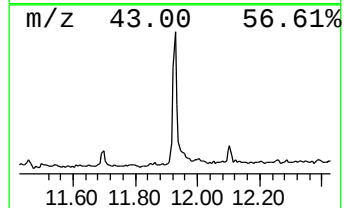
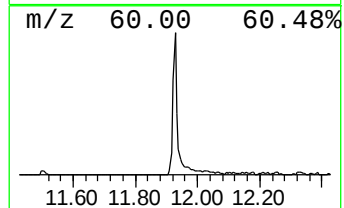
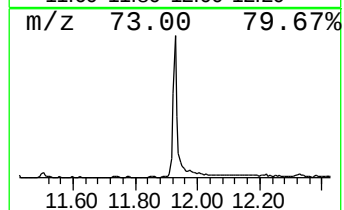
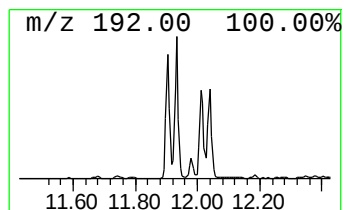
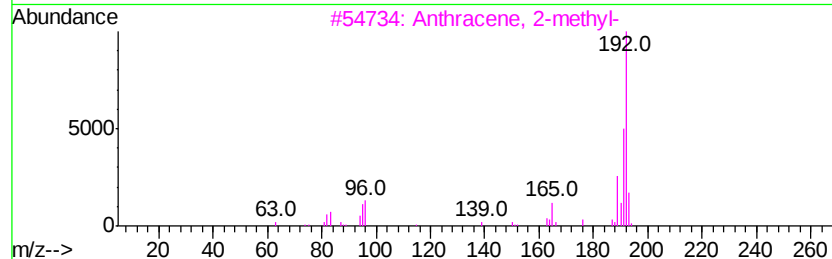
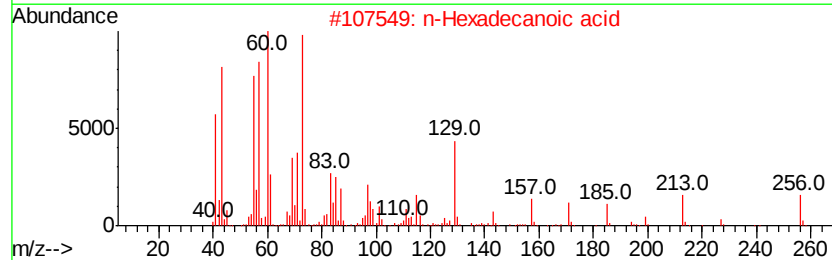
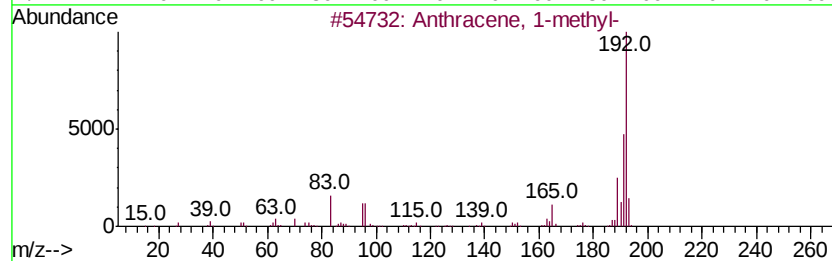
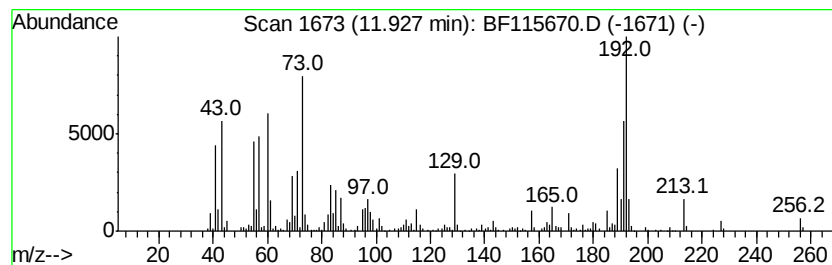
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Anthracene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	6.22 ng	530807	Phenanthrene-d10	11.40

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 1-methyl-	192	C15H12	000610-48-0	95
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071919\
Data File : BF115670.D
Acq On : 19 Jul 2019 12:07
Operator : HP/JU
Sample : K3618-05 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HD-01-071719

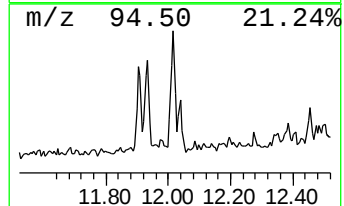
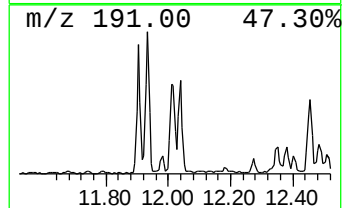
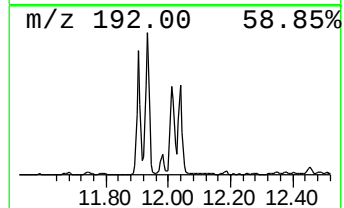
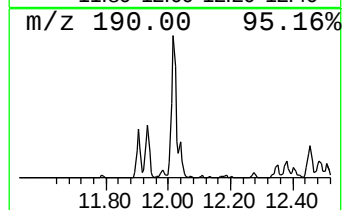
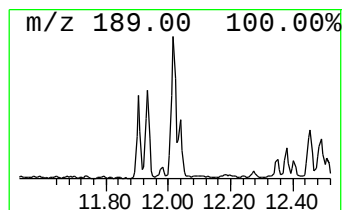
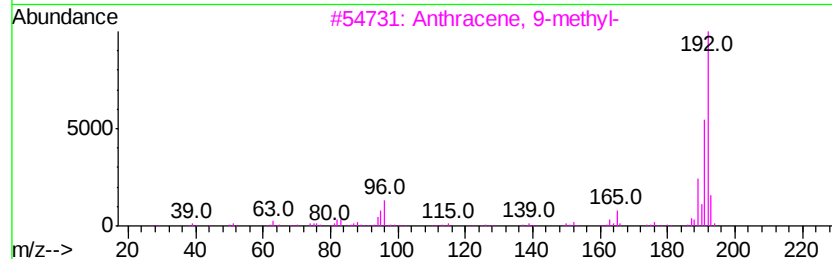
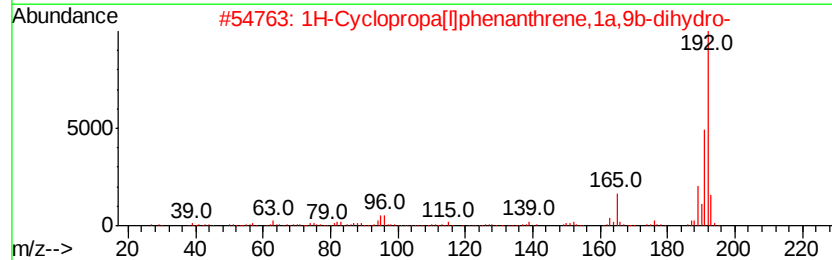
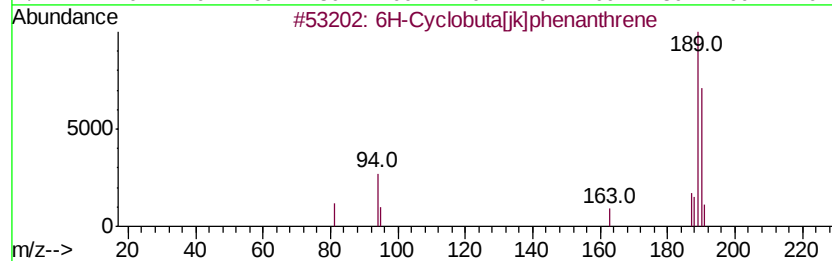
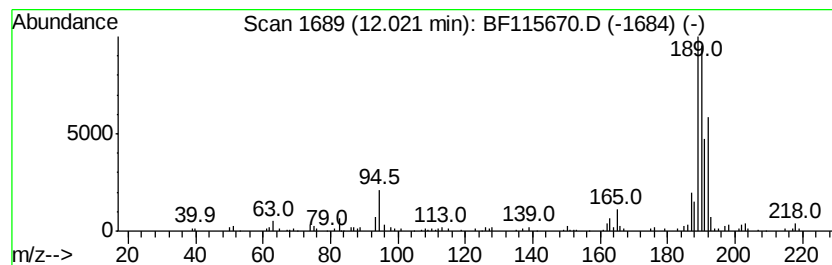
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown12.02 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.02	2.74 ng	234133	Phenanthrene-d10	11.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	49
2		1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	46
3		Anthracene, 9-methyl-	192	C15H12	000779-02-2	46
4		Anthracene, 1-methyl-	192	C15H12	000610-48-0	46
5		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	46



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071919\
Data File : BF115670.D
Acq On : 19 Jul 2019 12:07
Operator : HP/JU
Sample : K3618-05 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

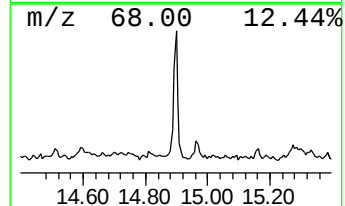
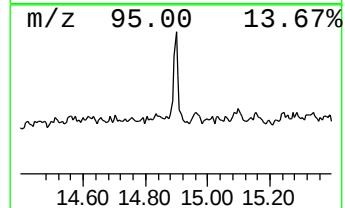
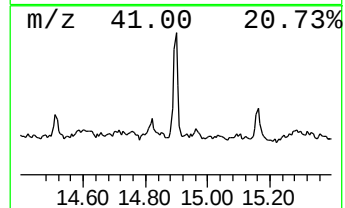
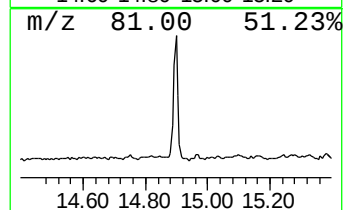
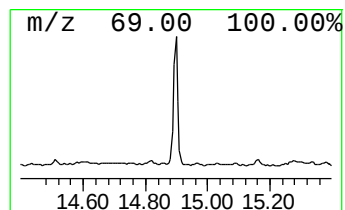
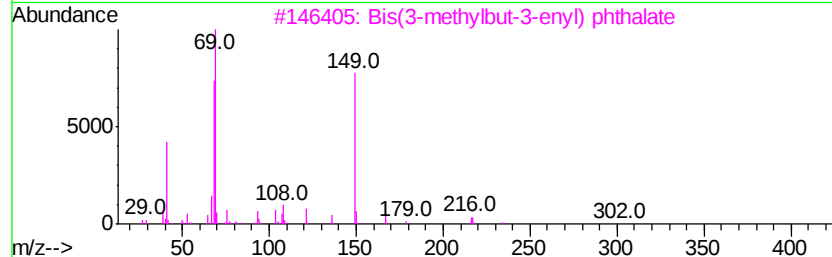
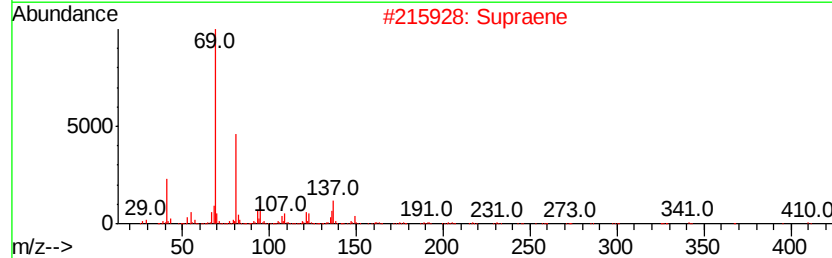
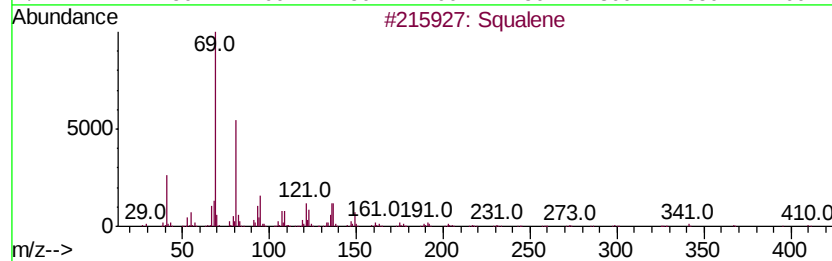
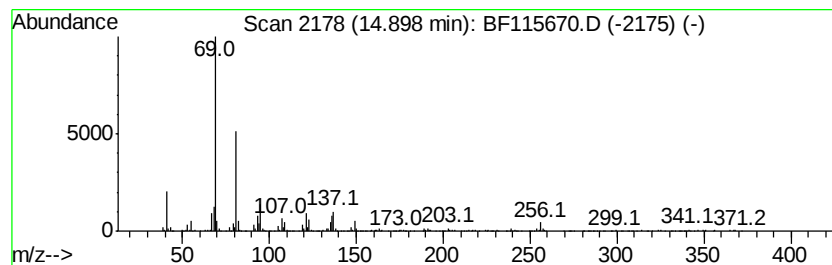
Instrument :
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ClientSampleId :
HD-01-071719

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Squalene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.90	4.78 ng	275471	Perylene-d12	15.52
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Squalene	410 C30H50	000111-02-4	87
2	Supraene	410 C30H50	007683-64-9	83
3	Bis(3-methylbut-3-enyl) phthalate	302 C18H22O4	022711-53-1	53
4	4-Methyl-1,5-Heptadiene	110 C8H14	000998-94-7	52
5	2,3,5-Trimethyl-2,3,5-hexanetric...	203 C12H17N3	003974-79-6	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071919\
Data File : BF115670.D
Acq On : 19 Jul 2019 12:07
Operator : HP/JU
Sample : K3618-05 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

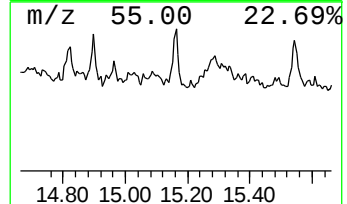
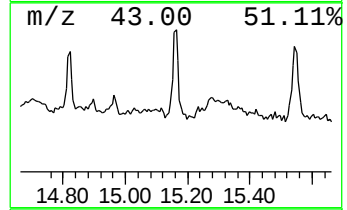
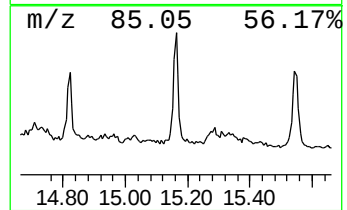
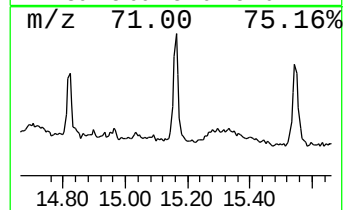
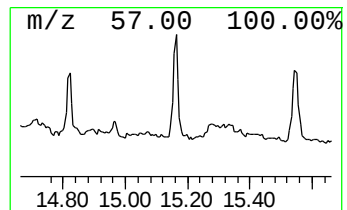
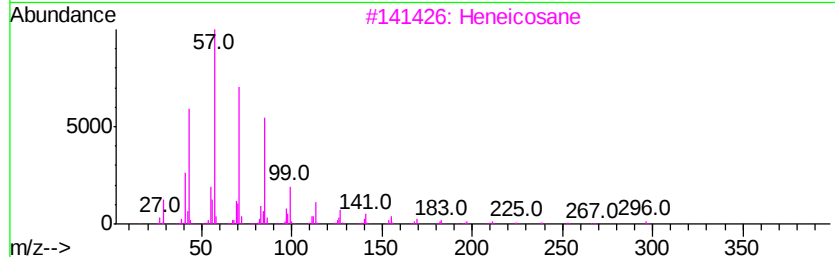
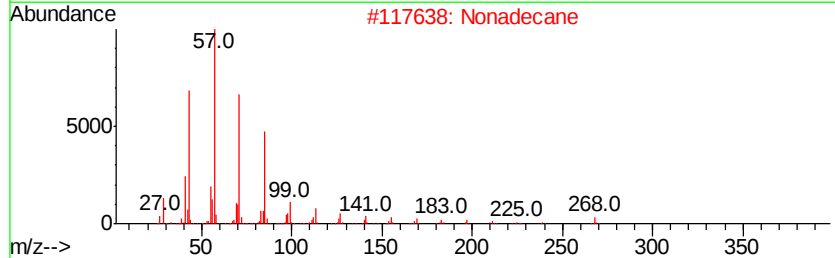
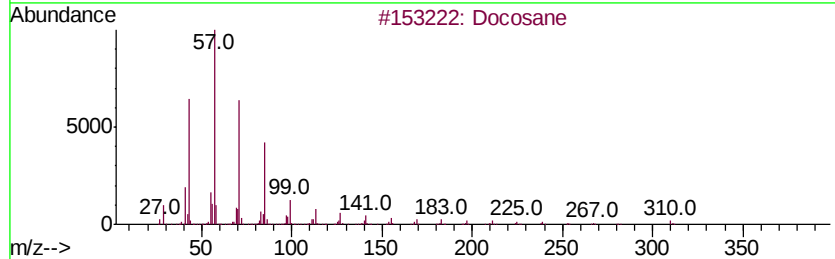
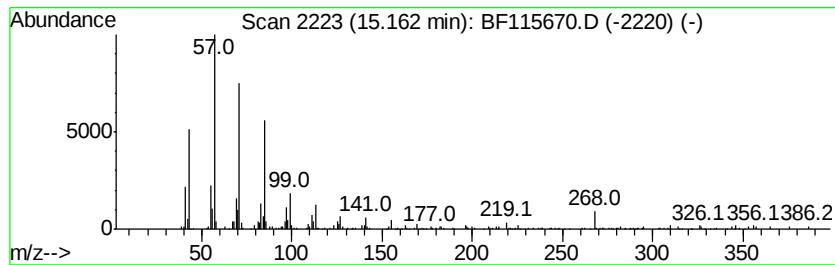
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HD-01-071719

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Docosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.16	2.68 ng	154401	Perylene-d12	15.52
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Docosane	310 C22H46	000629-97-0	95
2	Nonadecane	268 C19H40	000629-92-5	93
3	Heneicosane	296 C21H44	000629-94-7	90
4	Octacosane	394 C28H58	000630-02-4	90
5	Hexacosane	366 C26H54	000630-01-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071919\
Data File : BF115670.D
Acq On : 19 Jul 2019 12:07
Operator : HP/JU
Sample : K3618-05 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

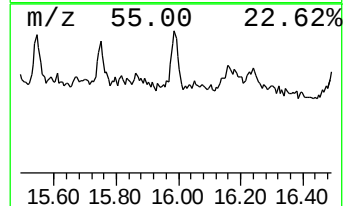
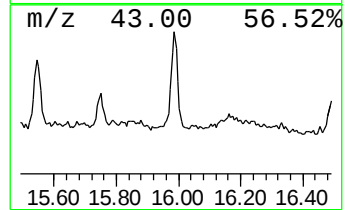
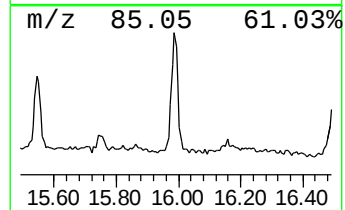
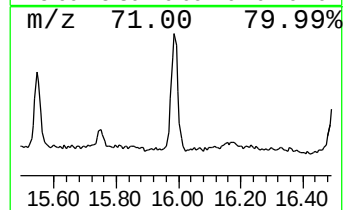
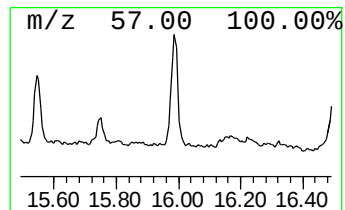
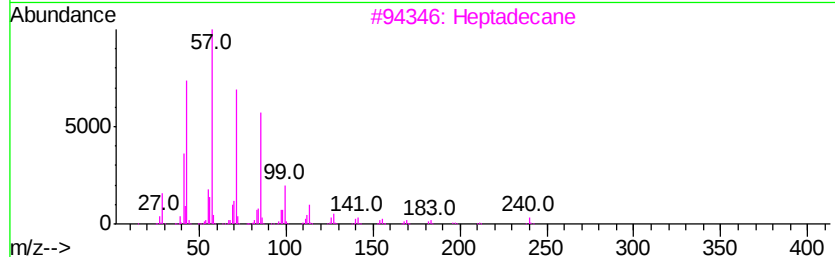
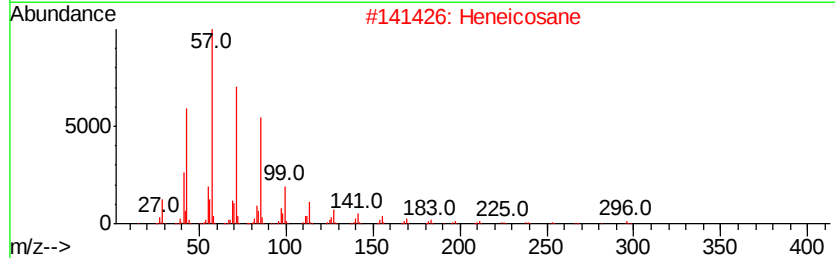
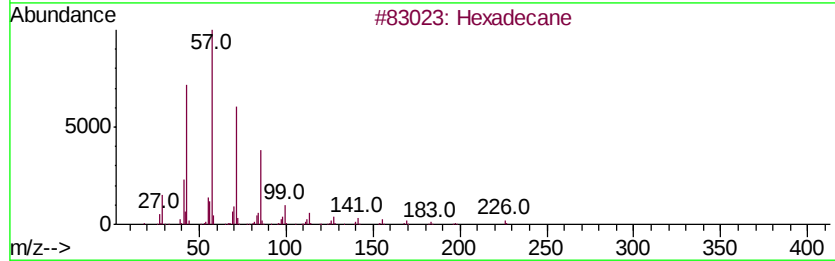
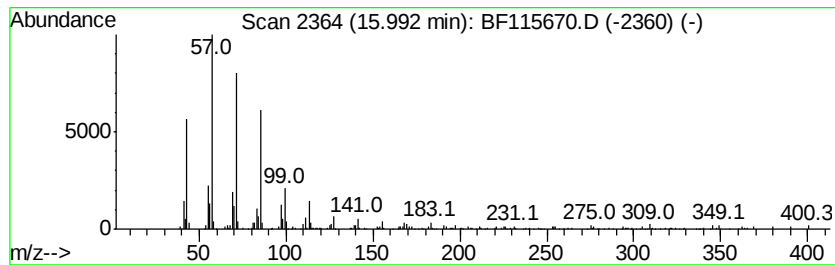
Instrument :
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ClientSampleId :
HD-01-071719

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.99	3.52 ng	202560	Perylene-d12	15.52
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexadecane	226 C16H34	000544-76-3	91
2	Heneicosane	296 C21H44	000629-94-7	91
3	Heptadecane	240 C17H36	000629-78-7	91
4	Octacosane	394 C28H58	000630-02-4	90
5	2-methyloctacosane	408 C29H60	1000376-72-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF071919\
Data File : BF115670.D
Acq On : 19 Jul 2019 12:07
Operator : HP/JU
Sample : K3618-05 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HD-01-071719

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF071219.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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unknown6.65	6.65	29.8	ng	1869290	1	6.88	1255110	20.0
Phenanthrene, 2-m...	11.90	2.4	ng	207917	4	11.40	1706590	20.0
Anthracene, 1-met...	11.93	6.2	ng	530807	4	11.40	1706590	20.0
unknown12.02	12.02	2.7	ng	234133	4	11.40	1706590	20.0
Squalene	14.90	4.8	ng	275471	6	15.52	1152380	20.0
Docosane	15.16	2.7	ng	154401	6	15.52	1152380	20.0
Hexadecane	15.99	3.5	ng	202560	6	15.52	1152380	20.0