

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
 Data File : BF117870.D
 Acq On : 19 Nov 2019 20:09
 Operator : JU
 Sample : K5865-12
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 3-(10-12)

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-BF111319.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	2.193	13	18	35	rVB	712238	1003726	25.42%	2.528%
2	4.510	408	412	419	rBV	375286	401488	10.17%	1.011%
3	5.122	508	516	520	rBV	1064380	1159443	29.36%	2.920%
4	5.522	580	584	587	rBV	1038210	881901	22.33%	2.221%
5	6.534	750	756	765	rBV	3178521	3217857	81.49%	8.105%
6	6.681	777	781	784	rBV	2279130	2013476	50.99%	5.071%
7	6.916	817	821	824	rBV	1493698	1245719	31.55%	3.138%
8	7.045	840	843	844	rBV	355295	262286	6.64%	0.661%
9	7.069	844	847	851	rVB	3072705	2766903	70.07%	6.969%
10	7.475	910	916	919	rBV	2507008	2111357	53.47%	5.318%
11	8.198	1036	1039	1042	rBV	1742671	1595168	40.40%	4.018%
12	9.280	1218	1223	1226	rBV	4867641	3948834	100.00%	9.946%
13	9.357	1233	1236	1239	rBV	79282	77636	1.97%	0.196%
14	9.622	1278	1281	1283	rBV	49582	48437	1.23%	0.122%
15	9.669	1287	1289	1294	rVB	464336	424047	10.74%	1.068%
16	9.892	1322	1327	1330	rVB	111464	94465	2.39%	0.238%
17	9.963	1333	1339	1343	rBV	2274358	1855354	46.98%	4.673%
18	10.163	1370	1373	1378	rBV3	61142	69726	1.77%	0.176%
19	10.204	1378	1380	1386	rVB3	55161	61788	1.56%	0.156%
20	10.280	1390	1393	1396	rBV	54022	51795	1.31%	0.130%
21	10.369	1404	1408	1410	rBV2	66678	75754	1.92%	0.191%
22	10.392	1410	1412	1415	rVB2	173906	129274	3.27%	0.326%
23	10.510	1428	1432	1434	rBV2	61298	64848	1.64%	0.163%
24	10.580	1439	1444	1446	rVV	69918	81490	2.06%	0.205%
25	10.616	1449	1450	1454	rVV2	67713	72917	1.85%	0.184%
26	10.669	1454	1459	1463	rVB5	48373	80788	2.05%	0.203%
27	10.751	1468	1473	1477	rVV2	203651	205000	5.19%	0.516%
28	10.792	1477	1480	1485	rVB3	37450	53633	1.36%	0.135%
29	10.869	1487	1493	1502	rVB2	190577	315322	7.99%	0.794%
30	10.957	1505	1508	1512	rBV2	44500	57199	1.45%	0.144%
31	11.063	1521	1526	1528	rBV3	87032	119811	3.03%	0.302%
32	11.092	1528	1531	1534	rVV	132088	138334	3.50%	0.348%
33	11.145	1538	1540	1543	rVB2	55668	49301	1.25%	0.124%
34	11.316	1563	1569	1572	rBV	192020	225746	5.72%	0.569%

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 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

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 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.351	1572	1575	1581	rVB2	109659	133814	3.39%	0.337%
36	11.457	1589	1593	1595	rBV	1994311	1695142	42.93%	4.269%
37	11.480	1595	1597	1600	rVV	651150	496792	12.58%	1.251%
38	11.527	1603	1605	1608	rVB	86033	80184	2.03%	0.202%
39	11.563	1608	1611	1614	rBV2	68904	80201	2.03%	0.202%
40	11.698	1630	1634	1639	rBV6	45485	83609	2.12%	0.211%
41	11.745	1639	1642	1645	rVB2	122089	100888	2.55%	0.254%
42	11.957	1675	1678	1680	rBV	255379	218035	5.52%	0.549%
43	11.986	1680	1683	1686	rVB	307277	261101	6.61%	0.658%
44	12.068	1694	1697	1699	rBV2	141287	136362	3.45%	0.343%
45	12.092	1699	1701	1705	rVB	142077	119976	3.04%	0.302%
46	12.157	1709	1712	1715	rBV	104706	87605	2.22%	0.221%
47	12.257	1726	1729	1731	rBV	68745	67746	1.72%	0.171%
48	12.404	1752	1754	1757	rBV	73294	54578	1.38%	0.137%
49	12.433	1757	1759	1762	rVV	65649	55219	1.40%	0.139%
50	12.510	1769	1772	1775	rBV	117234	119288	3.02%	0.300%
51	12.545	1775	1778	1781	rVV3	152546	156722	3.97%	0.395%
52	12.592	1784	1786	1790	rVV	63679	61088	1.55%	0.154%
53	12.674	1796	1800	1805	rBV	400430	378293	9.58%	0.953%
54	12.763	1813	1815	1818	rBV3	44975	52191	1.32%	0.131%
55	12.904	1834	1839	1840	rBV	463684	455214	11.53%	1.147%
56	12.921	1840	1842	1853	rVB	726965	938869	23.78%	2.365%
57	13.045	1859	1863	1870	rVB	3613739	3185924	80.68%	8.024%
58	13.127	1872	1877	1879	rBV3	51802	78032	1.98%	0.197%
59	13.227	1888	1894	1898	rVB2	72201	121188	3.07%	0.305%
60	13.274	1900	1902	1904	rBV2	57152	53043	1.34%	0.134%
61	13.333	1910	1912	1915	rVB2	60231	52520	1.33%	0.132%
62	13.621	1958	1961	1966	rBV2	50803	59340	1.50%	0.149%
63	13.739	1978	1981	1989	rVB	53237	60912	1.54%	0.153%
64	13.945	2013	2016	2019	rBV4	115035	164760	4.17%	0.415%
65	14.104	2037	2043	2046	rBV	1518694	1556057	39.41%	3.919%
66	14.127	2046	2047	2055	rVB	214032	139696	3.54%	0.352%
67	14.268	2068	2071	2075	rBV2	112121	130276	3.30%	0.328%
68	14.568	2119	2122	2126	rBV2	152948	163473	4.14%	0.412%
69	14.886	2173	2176	2180	rVB	186676	202837	5.14%	0.511%
70	14.968	2188	2190	2193	rBV	87184	84985	2.15%	0.214%
71	15.162	2219	2223	2231	rVB	149563	304130	7.70%	0.766%

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

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72	15.245	2233	2237	2241	rVV	196729	223295	5.65%	0.562%
73	15.474	2274	2276	2283	rVB	113344	156060	3.95%	0.393%
74	15.539	2284	2287	2291	rVV	104507	120880	3.06%	0.304%
75	15.609	2294	2299	2302	rVV	1011721	1318907	33.40%	3.322%
76	15.645	2302	2305	2313	rVB	261267	405661	10.27%	1.022%
77	16.104	2379	2383	2390	rVB	135229	206249	5.22%	0.519%
78	16.633	2470	2473	2480	rVB	72205	122201	3.09%	0.308%

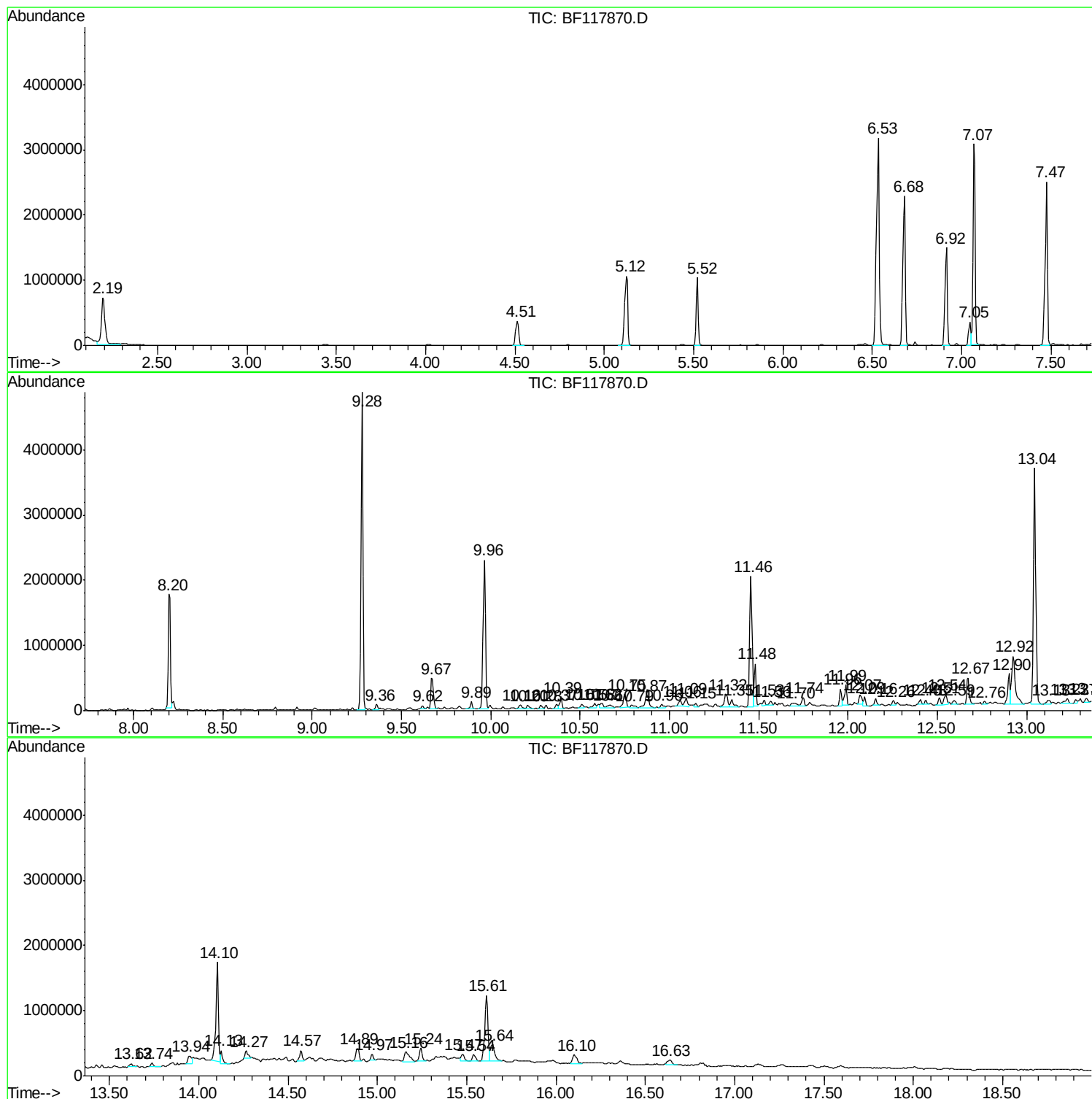
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Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
3-(10-12)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.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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Sample : K5865-12
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Instrument :
BNA_F
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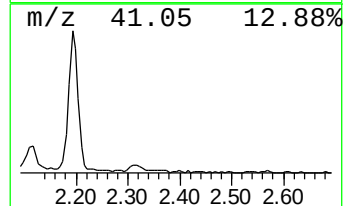
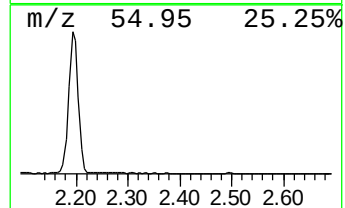
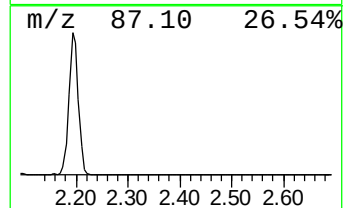
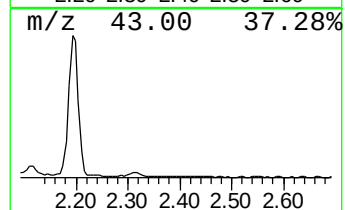
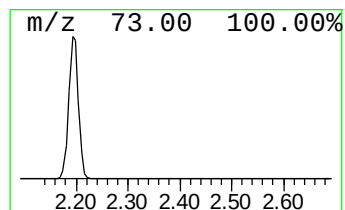
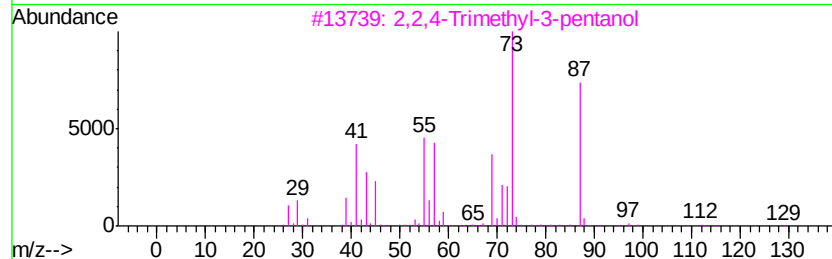
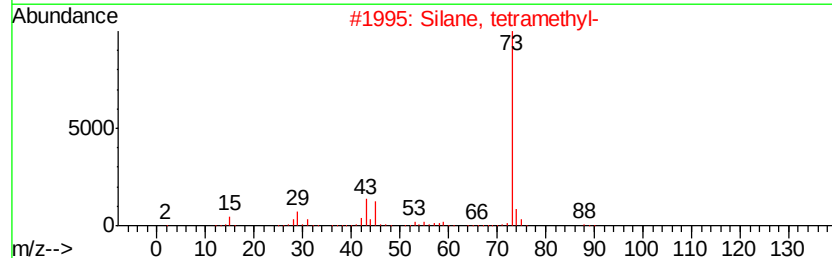
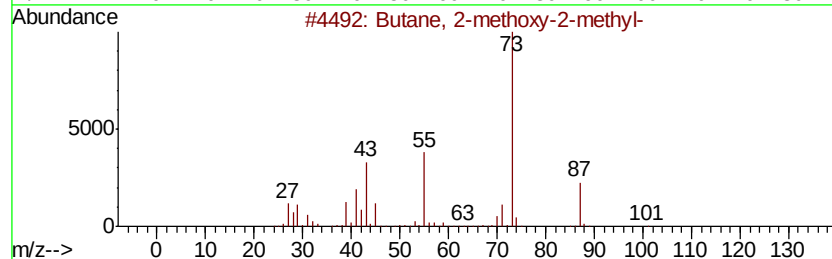
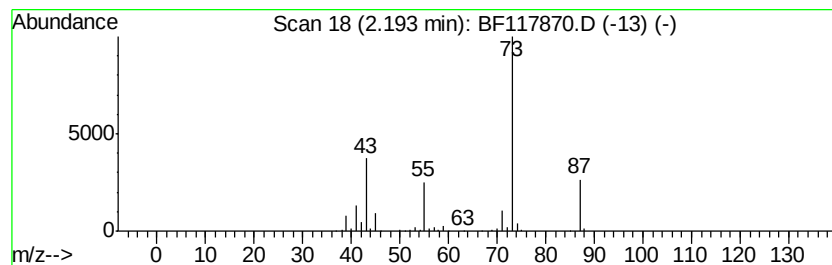
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
2.19	16.11 ng	1003730	1,4-Dichlorobenzene-d4	6.92

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Silane, tetramethyl-	88	C4H12Si	000075-76-3	43
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
5			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17



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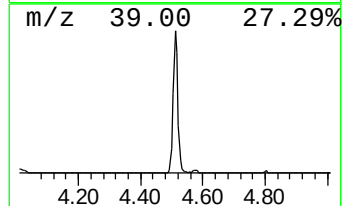
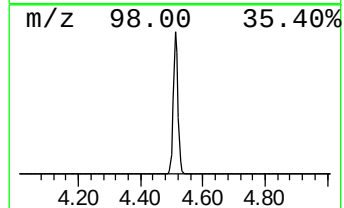
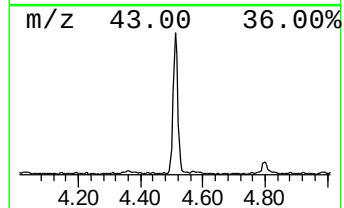
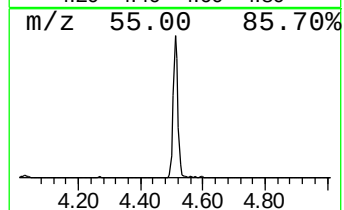
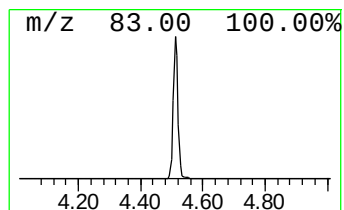
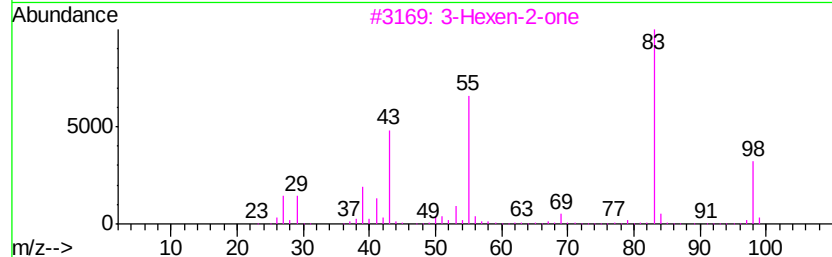
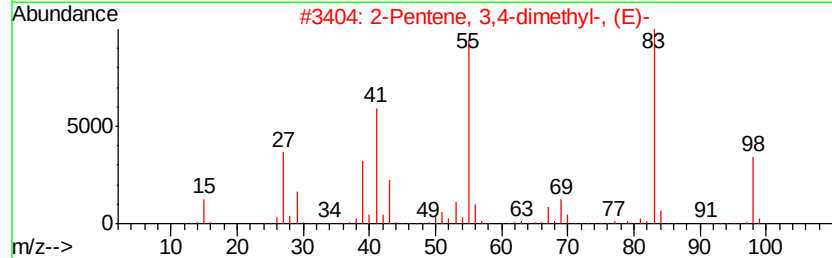
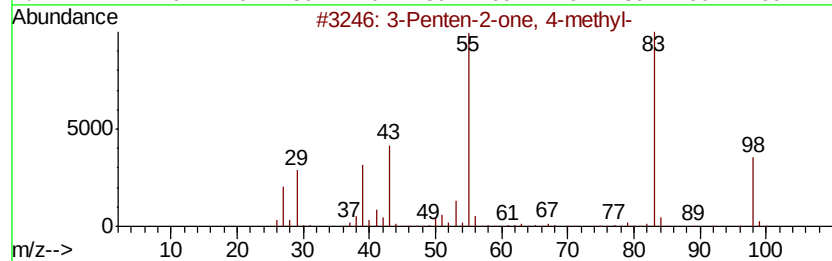
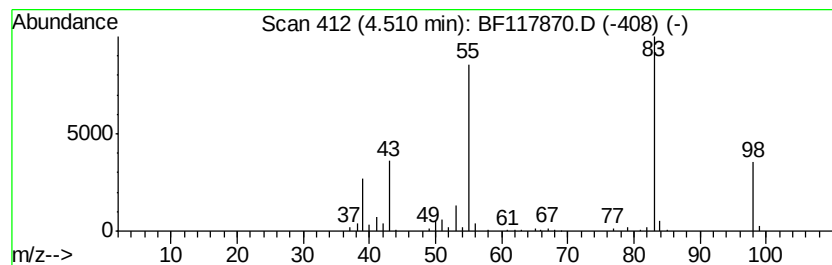
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 3-Penten-2-one, 4-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.51	6.45 ng	401488	1,4-Dichlorobenzene-d4	6.92

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	94
2			2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	90
3			3-Hexen-2-one	98	C6H10O	000763-93-9	87
4			2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	86
5			2-Pentene, 4,4-dimethyl-, (Z)-	98	C7H14	000762-63-0	80



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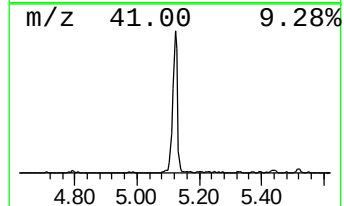
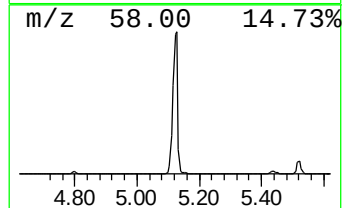
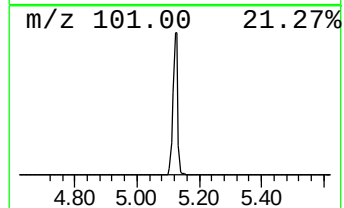
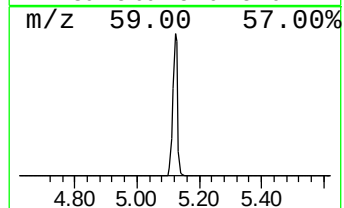
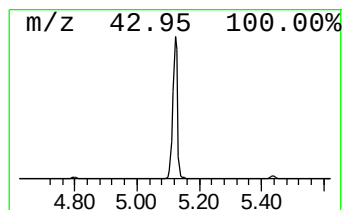
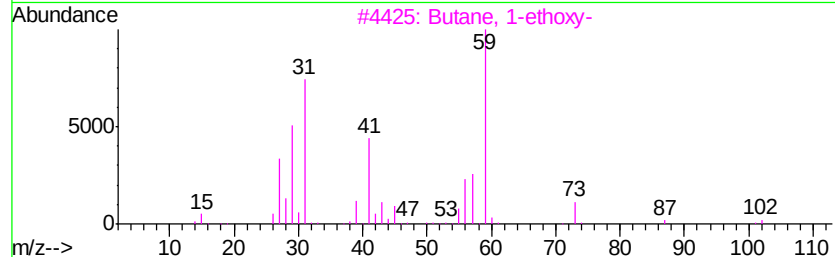
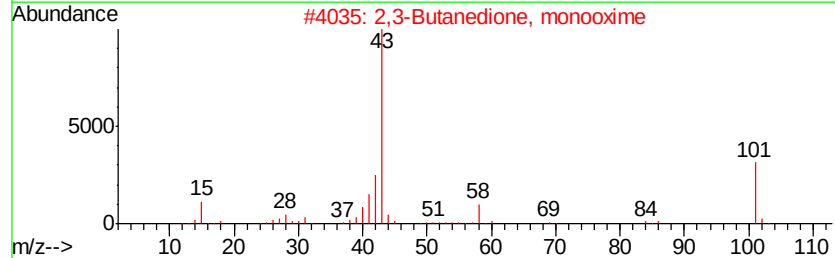
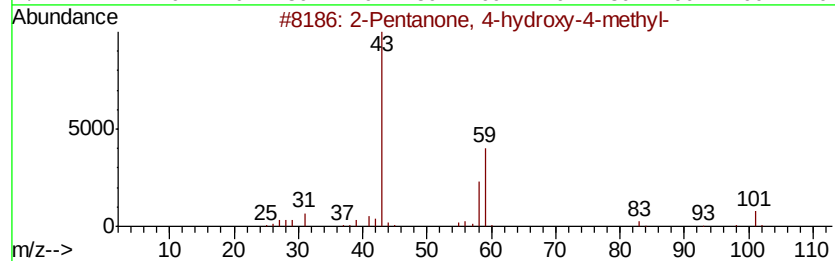
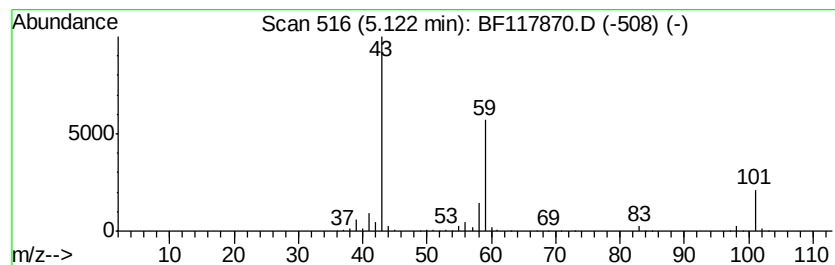
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	18.61 ng	1159440	1,4-Dichlorobenzene-d4	6.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
3		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9



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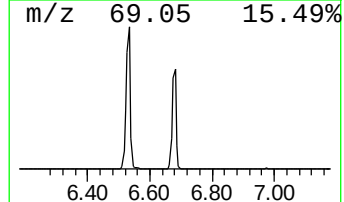
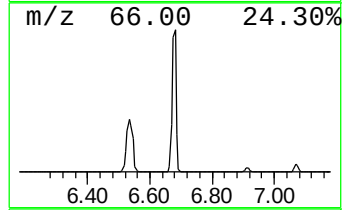
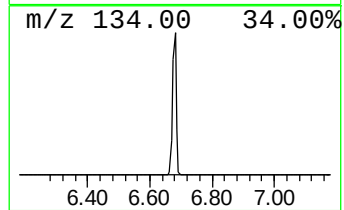
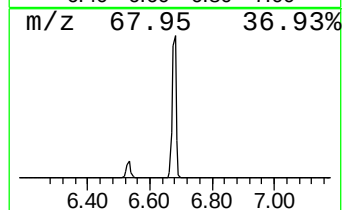
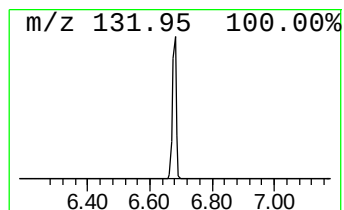
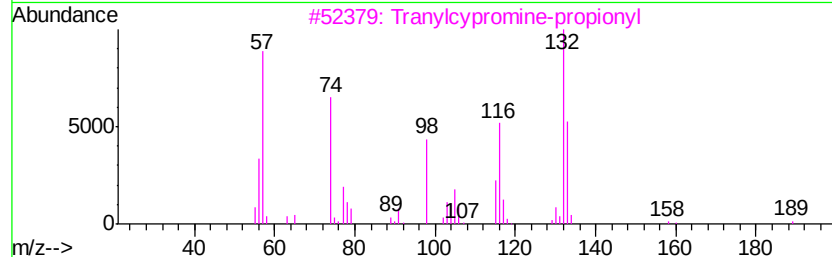
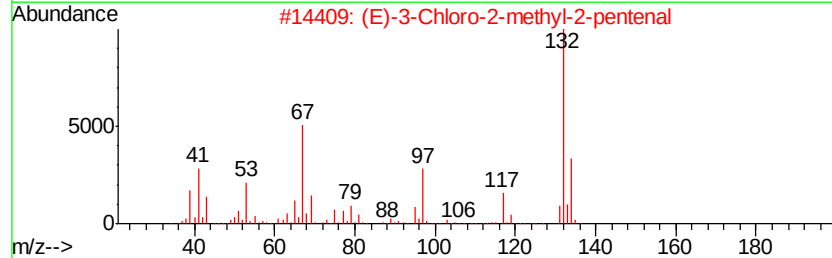
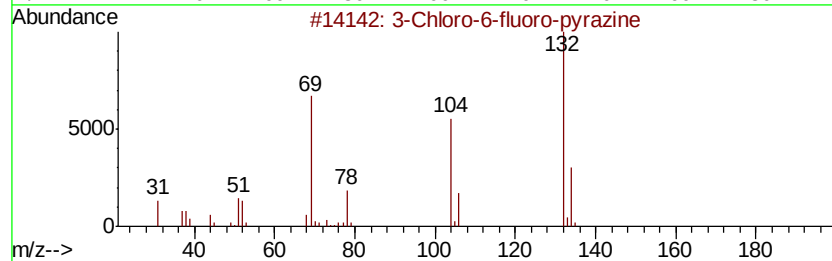
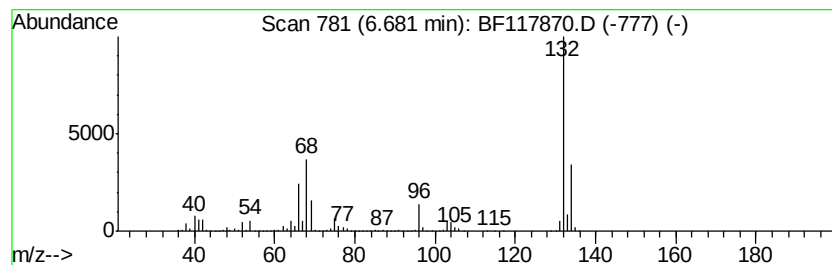
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 unknown6.68 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.68	32.33 ng	2013480	1,4-Dichlorobenzene-d4	6.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	16
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
Data File : BF117870.D
Acq On : 19 Nov 2019 20:09
Operator : JU
Sample : K5865-12
Misc :
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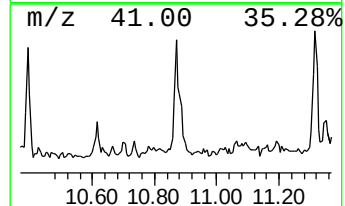
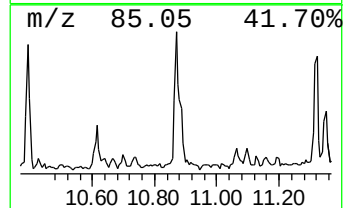
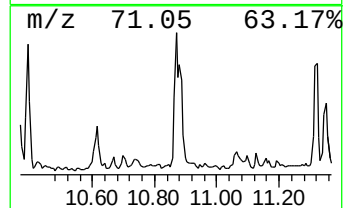
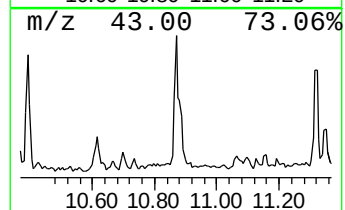
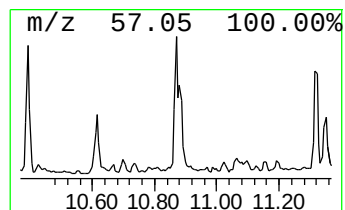
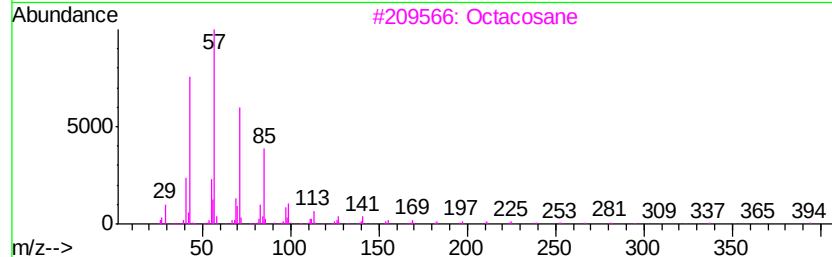
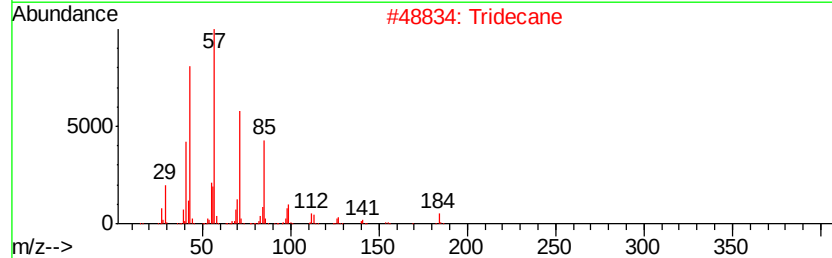
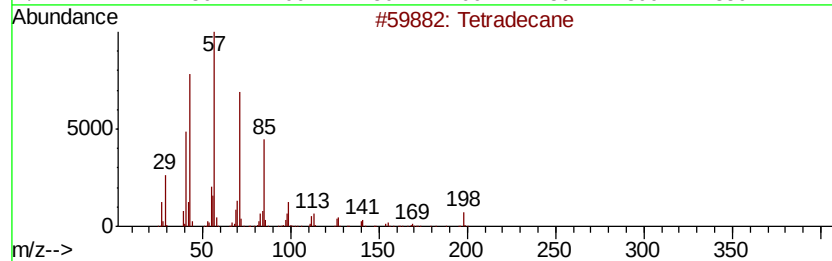
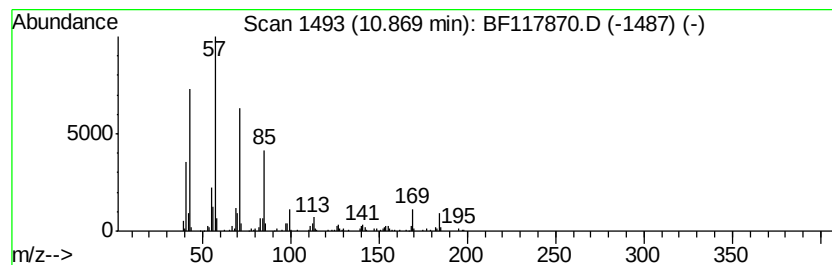
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Tetradecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.87	3.72 ng	315322	Phenanthrene-d10		11.46
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	93
2	Tridecane	184	C13H28	000629-50-5	87
3	Octacosane	394	C28H58	000630-02-4	74
4	Tridecane, 7-hexyl-	268	C19H40	007225-66-3	74
5	Eicosane	282	C20H42	000112-95-8	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
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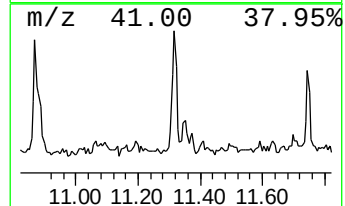
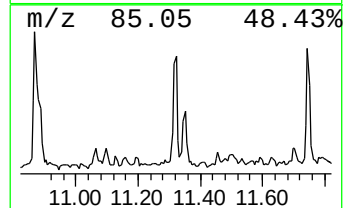
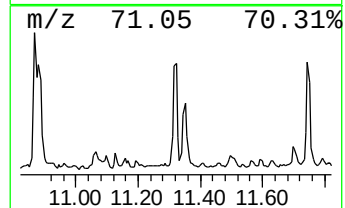
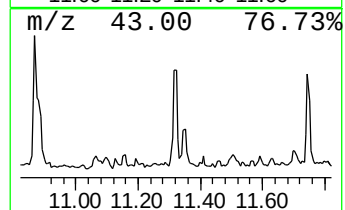
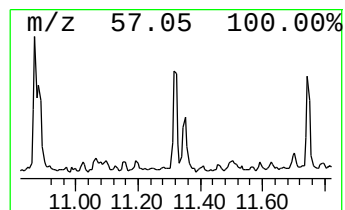
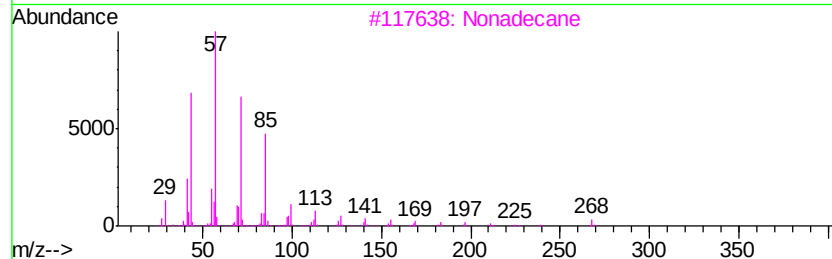
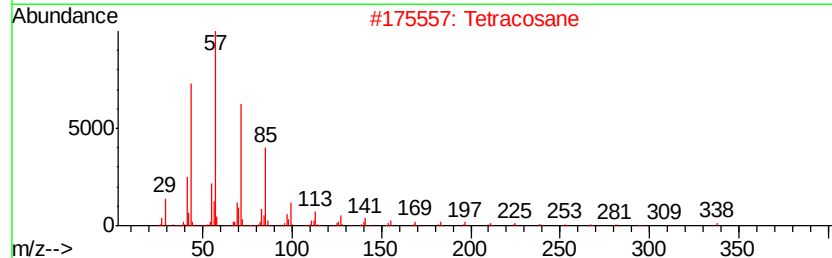
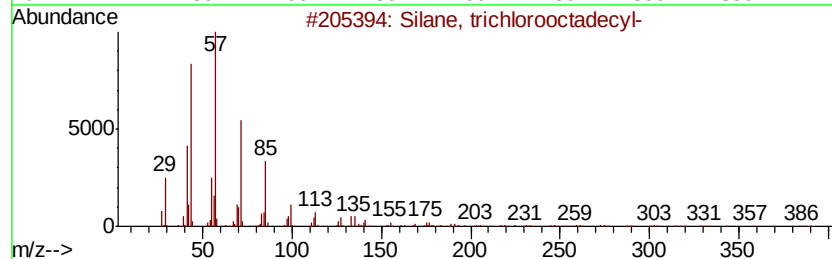
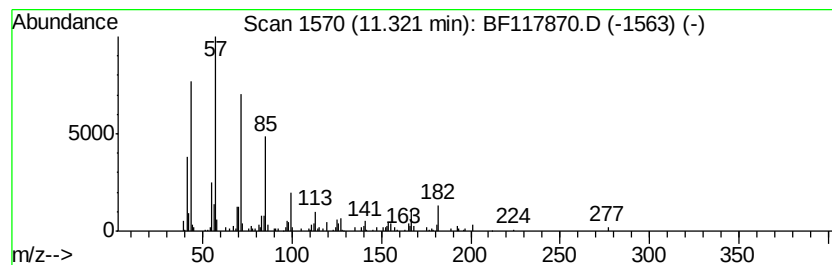
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Silane, trichlorooctadecyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.32	2.66 ng	225746	Phenanthrene-d10		11.46	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Silane, trichlorooctadecyl-		386	C18H37Cl3Si	000112-04-9	53
2	Tetracosane		338	C24H50	000646-31-1	53
3	Nonadecane		268	C19H40	000629-92-5	53
4	Tetradecane		198	C14H30	000629-59-4	53
5	Heptadecane		240	C17H36	000629-78-7	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
Data File : BF117870.D
Acq On : 19 Nov 2019 20:09
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Sample : K5865-12
Misc :
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Instrument :
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ClientSampleId :
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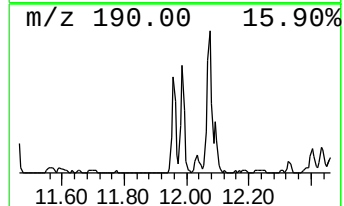
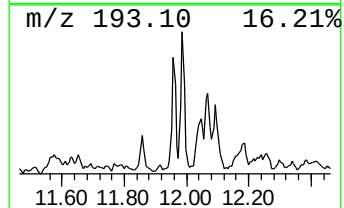
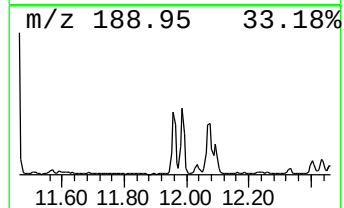
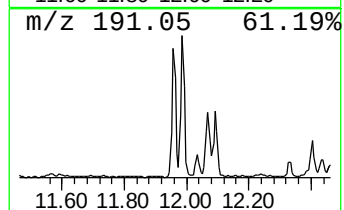
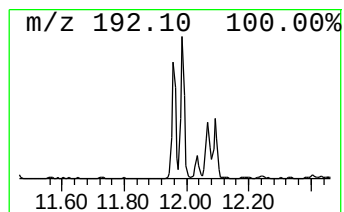
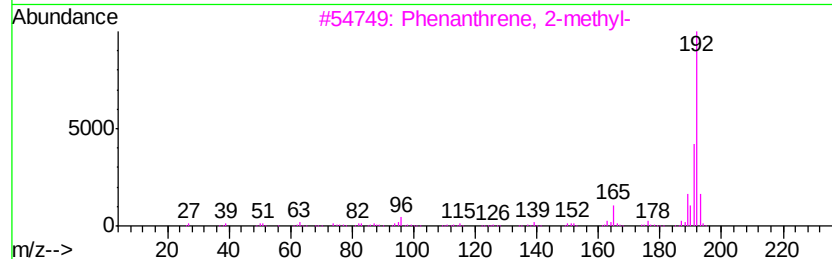
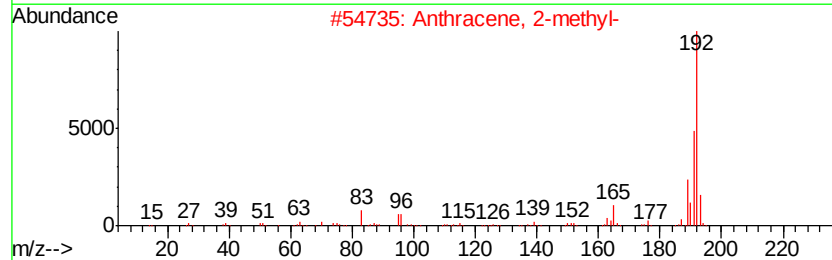
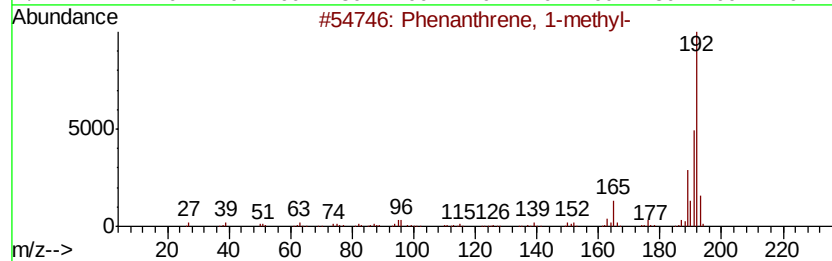
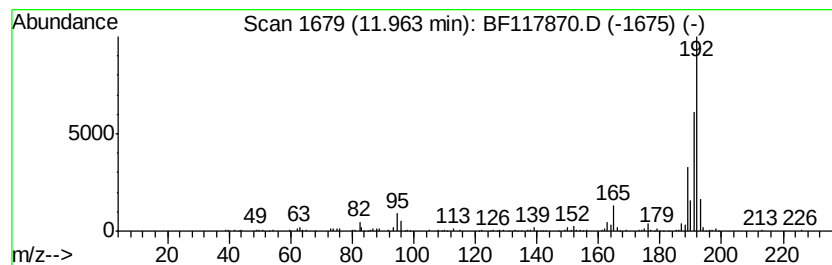
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Phenanthrene, 1-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.96	2.57 ng	218035	Phenanthrene-d10	11.46

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
4		Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
5		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
Data File : BF117870.D
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Misc :
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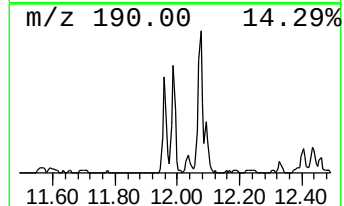
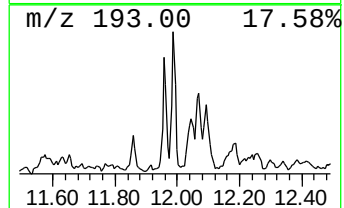
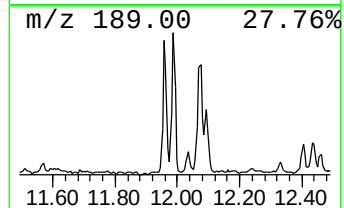
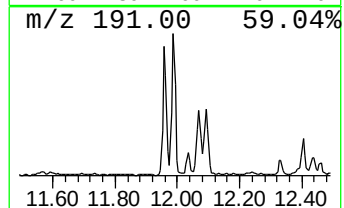
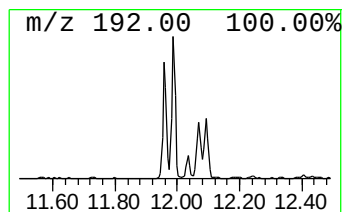
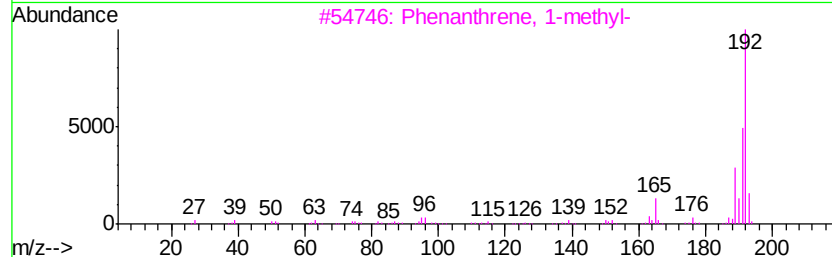
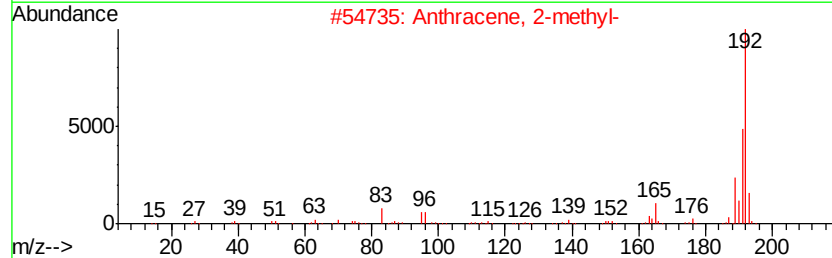
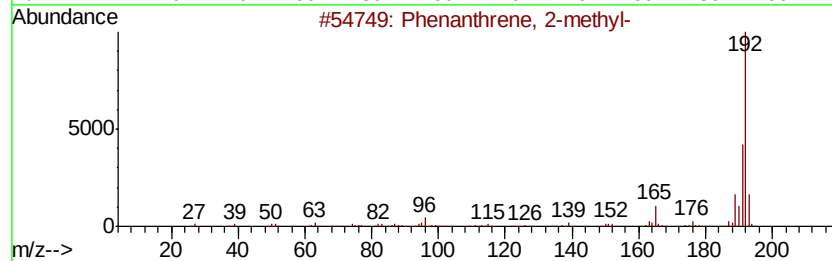
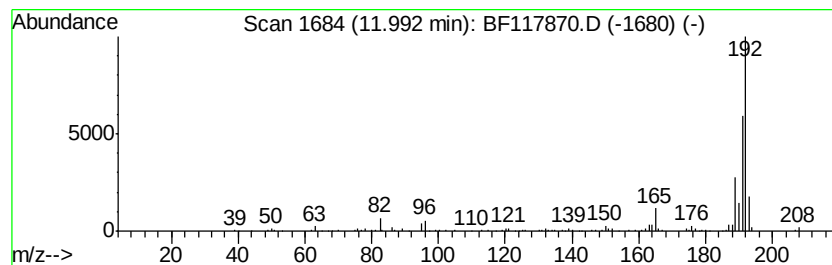
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Phenanthrene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.99	3.08 ng	261101	Phenanthrene-d10	11.46

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	97
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
4		1H-Cyclopropa[1,1]phenanthrene, 1a, ...	192	C15H12	000949-41-7	96
5		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
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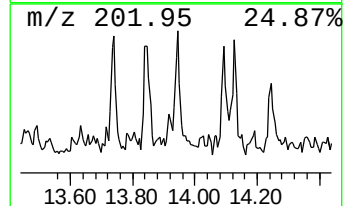
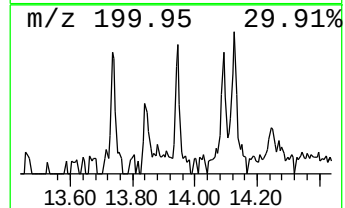
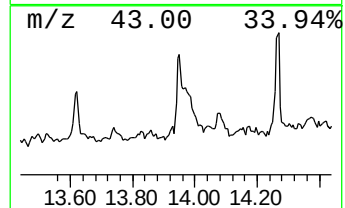
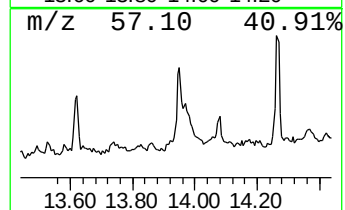
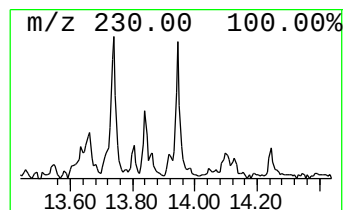
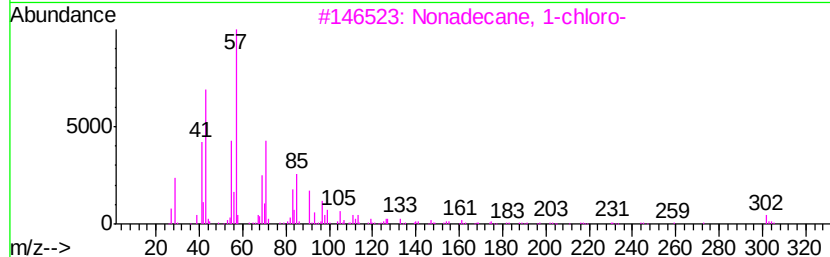
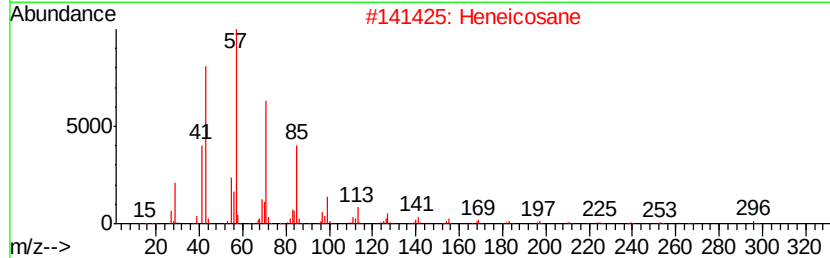
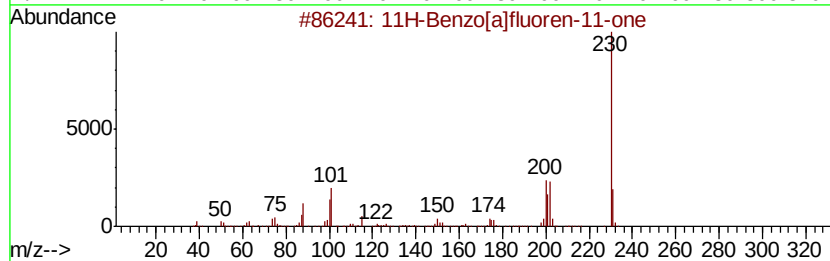
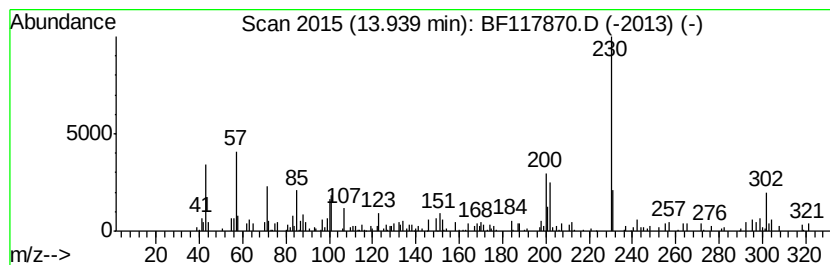
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 11H-Benzo[a]fluoren-11-one Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.94	2.12 ng	164760	Chrysene-d12	14.10	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	91
2	Heneicosane	296	C21H44	000629-94-7	56
3	Nonadecane, 1-chloro-	302	C19H39Cl	062016-76-6	47
4	Nonadecane	268	C19H40	000629-92-5	46
5	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	45



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
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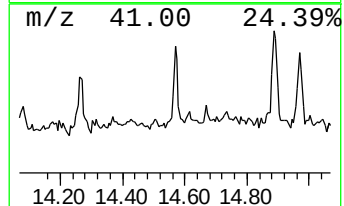
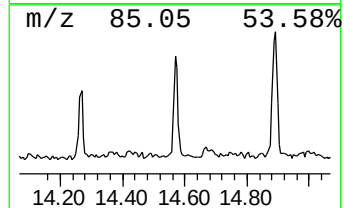
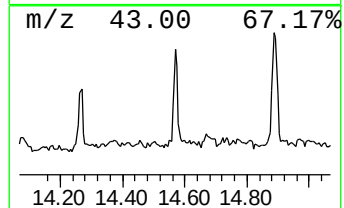
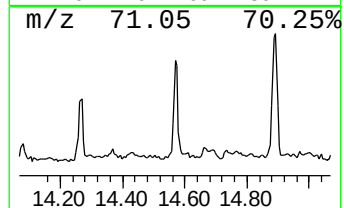
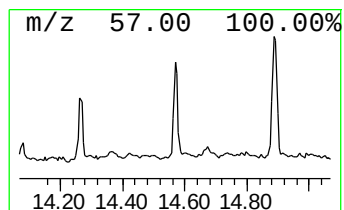
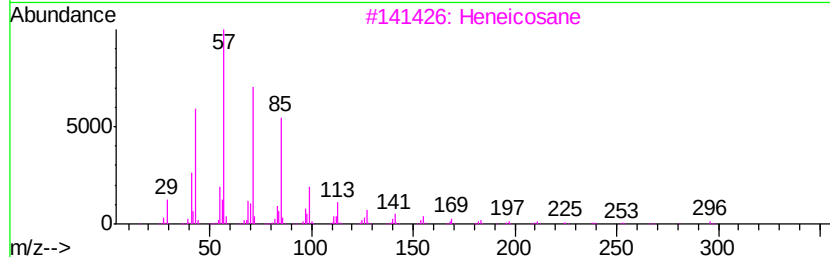
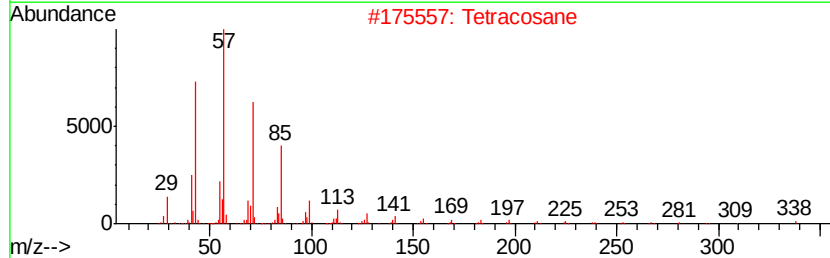
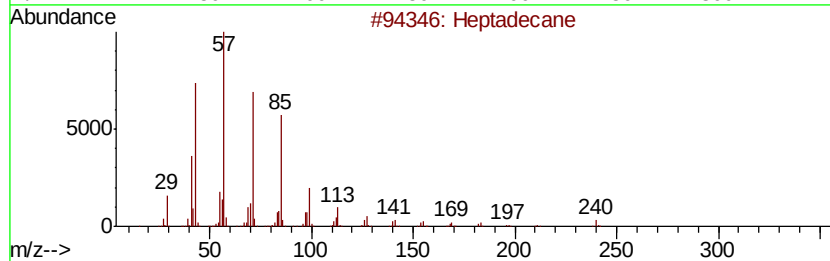
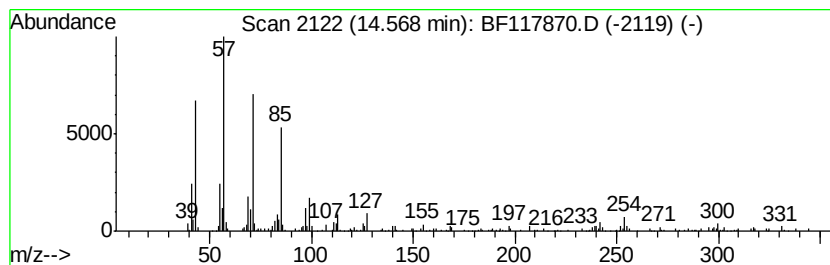
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heptadecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.57	2.10 ng	163473	Chrysene-d12	14.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	93
2		Tetracosane	338	C24H50	000646-31-1	93
3		Heneicosane	296	C21H44	000629-94-7	83
4		Hexadecane	226	C16H34	000544-76-3	81
5		Octadecane	254	C18H38	000593-45-3	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
Data File : BF117870.D
Acq On : 19 Nov 2019 20:09
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ClientSampleId :
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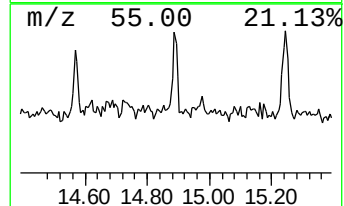
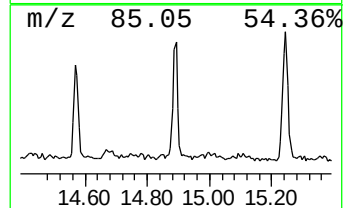
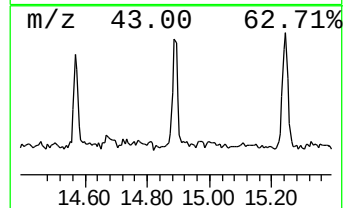
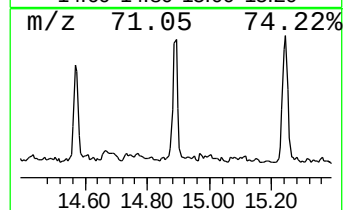
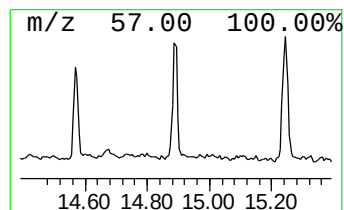
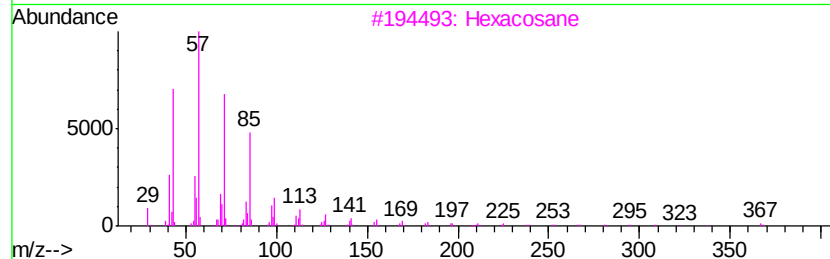
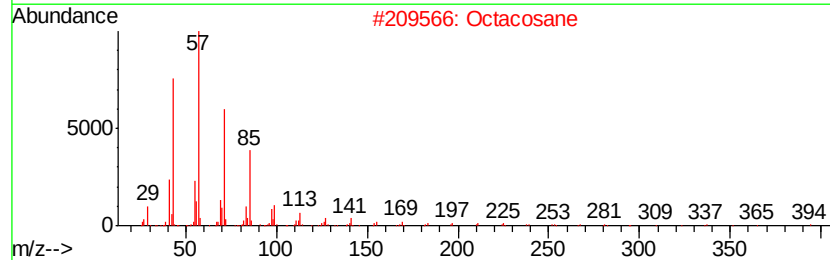
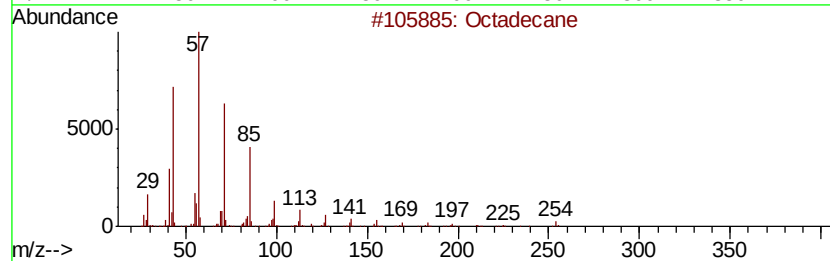
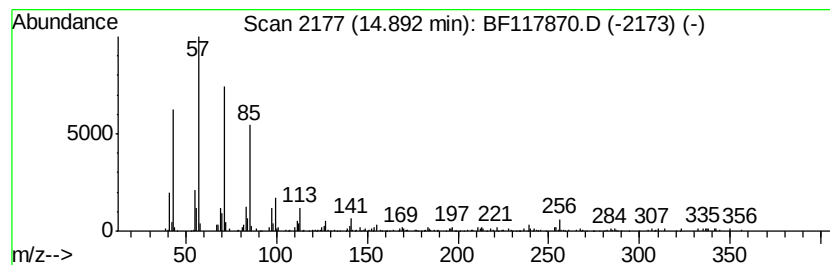
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Octadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.89	3.08 ng	202837	Perylene-d12	15.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	91
2		Octacosane	394	C28H58	000630-02-4	90
3		Hexacosane	366	C26H54	000630-01-3	90
4		Pentadecane	212	C15H32	000629-62-9	87
5		Heptadecane	240	C17H36	000629-78-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
Data File : BF117870.D
Acq On : 19 Nov 2019 20:09
Operator : JU
Sample : K5865-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
3-(10-12)

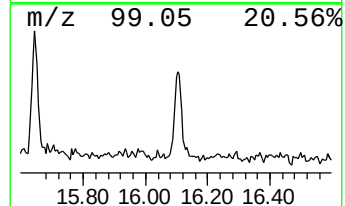
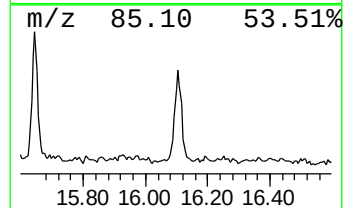
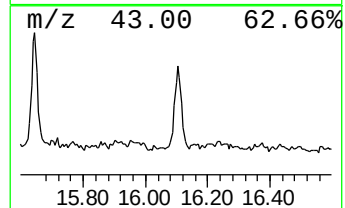
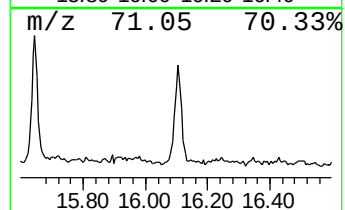
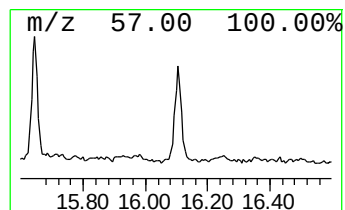
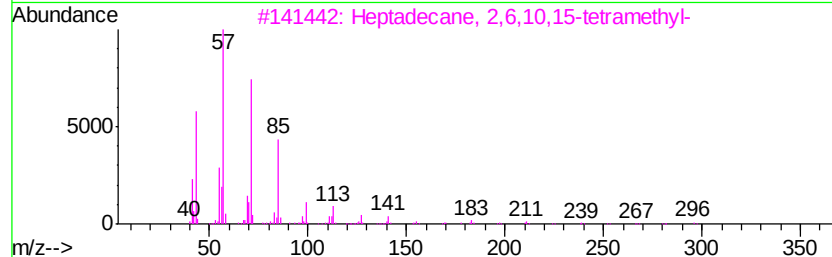
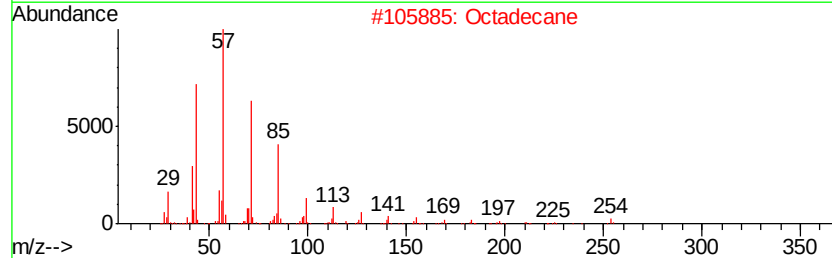
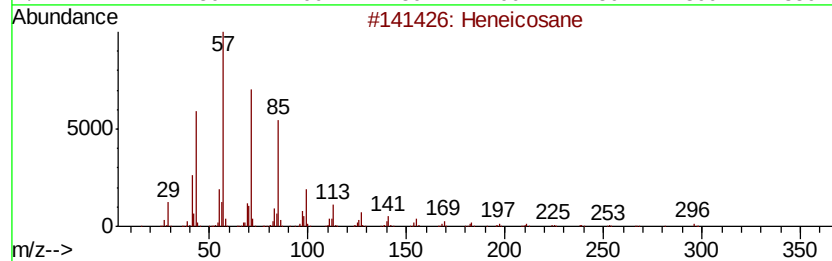
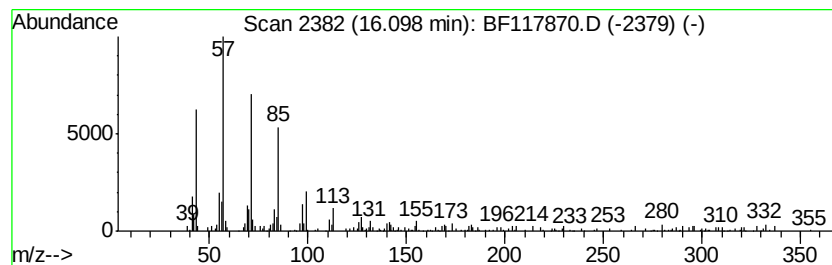
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Heneicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.10	3.13 ng	206249	Perylene-d12	15.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	97
2		Octadecane	254	C18H38	000593-45-3	91
3		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
4		Hexacosane	366	C26H54	000630-01-3	90
5		2-methyloctacosane	408	C29H60	1000376-72-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF111919\
Data File : BF117870.D
Acq On : 19 Nov 2019 20:09
Operator : JU
Sample : K5865-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
3-(10-12)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.19	16.1	ng	1003730	1	6.92	1245720	20.0
3-Penten-2-one, 4...	4.51	6.5	ng	401488	1	6.92	1245720	20.0
2-Pentanone, 4-hy...	5.12	18.6	ng	1159440	1	6.92	1245720	20.0
unknown6.68	6.68	32.3	ng	2013480	1	6.92	1245720	20.0
Tetradecane	10.87	3.7	ng	315322	4	11.46	1695140	20.0
Silane, trichloro...	11.32	2.7	ng	225746	4	11.46	1695140	20.0
Phenanthrene, 1-m...	11.96	2.6	ng	218035	4	11.46	1695140	20.0
Phenanthrene, 2-m...	11.99	3.1	ng	261101	4	11.46	1695140	20.0
11H-Benzo[a]fluor...	13.94	2.1	ng	164760	5	14.10	1556060	20.0
Heptadecane	14.57	2.1	ng	163473	5	14.10	1556060	20.0
Octadecane	14.89	3.1	ng	202837	6	15.61	1318910	20.0
Heneicosane	16.10	3.1	ng	206249	6	15.61	1318910	20.0