

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF081319\
 Data File : BF116233.D
 Acq On : 13 Aug 2019 15:54
 Operator : HP/JU
 Sample : K4227-21
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 1S-I(0-1)-COMP

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-BF073019.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	2.134	3	8	30	rVB	1511403	2117236	48.24%	4.237%
2	5.463	570	574	604	rBV	2814632	3080810	70.20%	6.165%
3	6.475	741	746	764	rBV	2856514	3427664	78.11%	6.860%
4	6.610	764	769	772	rBV	3108736	3165345	72.13%	6.335%
5	6.834	803	807	818	rVB	963576	876311	19.97%	1.754%
6	6.987	829	833	849	rBV	2557880	2616671	59.63%	5.237%
7	7.398	899	903	933	rBV	2103978	2236639	50.97%	4.476%
8	8.110	1021	1024	1042	rBV	1013676	1181160	26.91%	2.364%
9	9.186	1202	1207	1220	rBV	3556385	3919445	89.31%	7.844%
10	9.586	1271	1275	1285	rBV	318877	376428	8.58%	0.753%
11	9.869	1319	1323	1327	rBV	1611705	1435233	32.70%	2.872%
12	9.898	1327	1328	1337	rVB	74283	70216	1.60%	0.141%
13	10.422	1414	1417	1423	rBV	40922	44089	1.00%	0.088%
14	10.663	1454	1458	1472	rBV	2407960	2693332	61.37%	5.390%
15	11.257	1554	1559	1563	rVB2	45726	56668	1.29%	0.113%
16	11.357	1572	1576	1578	rBV	1698880	1546878	35.25%	3.096%
17	11.380	1578	1580	1585	rVV	907064	827281	18.85%	1.656%
18	11.439	1585	1590	1593	rVB	181328	256696	5.85%	0.514%
19	11.598	1613	1617	1622	rBV	58933	88706	2.02%	0.178%
20	11.863	1659	1662	1663	rBV	113438	106619	2.43%	0.213%
21	11.886	1663	1666	1672	rVV2	449647	583697	13.30%	1.168%
22	11.939	1672	1675	1678	rVV	81929	101509	2.31%	0.203%
23	11.974	1678	1681	1683	rVV	227161	246387	5.61%	0.493%
24	11.992	1683	1684	1690	rVB	93762	91600	2.09%	0.183%
25	12.151	1708	1711	1716	rBV	93241	98578	2.25%	0.197%
26	12.198	1716	1719	1722	rBV2	58967	50531	1.15%	0.101%
27	12.304	1732	1737	1740	rBV	48762	62962	1.43%	0.126%
28	12.410	1751	1755	1757	rBV	118375	112777	2.57%	0.226%
29	12.504	1767	1771	1775	rBV2	57624	83619	1.91%	0.167%
30	12.574	1779	1783	1788	rBV	1689605	1623331	36.99%	3.249%
31	12.663	1796	1798	1805	rBV3	50436	68986	1.57%	0.138%
32	12.727	1805	1809	1811	rVB2	69538	63496	1.45%	0.127%
33	12.804	1815	1822	1828	rBV	1392187	1659763	37.82%	3.322%
34	12.857	1828	1831	1835	rVB2	88635	97462	2.22%	0.195%

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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-BF073019.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	12.945	1838	1846	1849	rBV	3861030	4388529	100.00%	8.783%
36	12.974	1849	1851	1855	rVB	69328	75399	1.72%	0.151%
37	13.016	1855	1858	1863	rBV3	88987	163860	3.73%	0.328%
38	13.110	1870	1874	1875	rBV2	63417	79437	1.81%	0.159%
39	13.133	1875	1878	1881	rVB	205901	185750	4.23%	0.372%
40	13.198	1886	1889	1892	rVV	155264	140195	3.19%	0.281%
41	13.233	1892	1895	1902	rVV3	125825	208431	4.75%	0.417%
42	13.327	1908	1911	1914	rBV	70031	73695	1.68%	0.147%
43	13.363	1914	1917	1923	rBV2	74603	100961	2.30%	0.202%
44	13.616	1957	1960	1963	rBV3	44434	59495	1.36%	0.119%
45	13.651	1963	1966	1969	rVB	88237	84412	1.92%	0.169%
46	13.757	1979	1984	1985	rBV2	138688	150155	3.42%	0.300%
47	13.774	1985	1987	1990	rVB	124568	97583	2.22%	0.195%
48	13.804	1990	1992	1994	rBV	91352	95552	2.18%	0.191%
49	13.833	1994	1997	2000	rBV2	291778	437735	9.97%	0.876%
50	14.004	2017	2026	2028	rBV2	1711431	2385286	54.35%	4.774%
51	14.027	2028	2030	2037	rVB	759567	698256	15.91%	1.397%
52	14.110	2041	2044	2047	rVB	173356	153702	3.50%	0.308%
53	14.145	2047	2050	2052	rBV2	54792	56195	1.28%	0.112%
54	14.392	2087	2092	2095	rVV	134472	162170	3.70%	0.325%
55	14.451	2099	2102	2103	rVV2	73325	69403	1.58%	0.139%
56	14.492	2106	2109	2111	rVV2	53472	48293	1.10%	0.097%
57	14.721	2144	2148	2150	rBV3	63516	90584	2.06%	0.181%
58	14.751	2150	2153	2156	rVB	116188	128021	2.92%	0.256%
59	14.827	2161	2166	2170	rVB	361953	396992	9.05%	0.794%
60	14.892	2173	2177	2180	rBV2	69721	86122	1.96%	0.172%
61	15.045	2199	2203	2206	rBV	529057	795999	18.14%	1.593%
62	15.157	2218	2222	2228	rVB	106215	128072	2.92%	0.256%
63	15.345	2250	2254	2261	rVB	274670	355026	8.09%	0.710%
64	15.410	2261	2265	2269	rBV	428795	526812	12.00%	1.054%
65	15.468	2269	2275	2280	rBV	846606	1217556	27.74%	2.437%
66	15.880	2340	2345	2350	rBV	124241	169450	3.86%	0.339%
67	15.974	2355	2361	2362	rBV4	37703	66521	1.52%	0.133%
68	16.039	2370	2372	2383	rVB2	36730	61185	1.39%	0.122%
69	16.374	2422	2429	2442	rVB3	80892	176569	4.02%	0.353%
70	16.951	2520	2527	2537	rBV	186349	397988	9.07%	0.796%
71	17.127	2551	2557	2563	rBV3	51056	121411	2.77%	0.243%

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073019.M
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72	17.398	2597	2603	2618	rVB2	107455	267700	6.10%	0.536%
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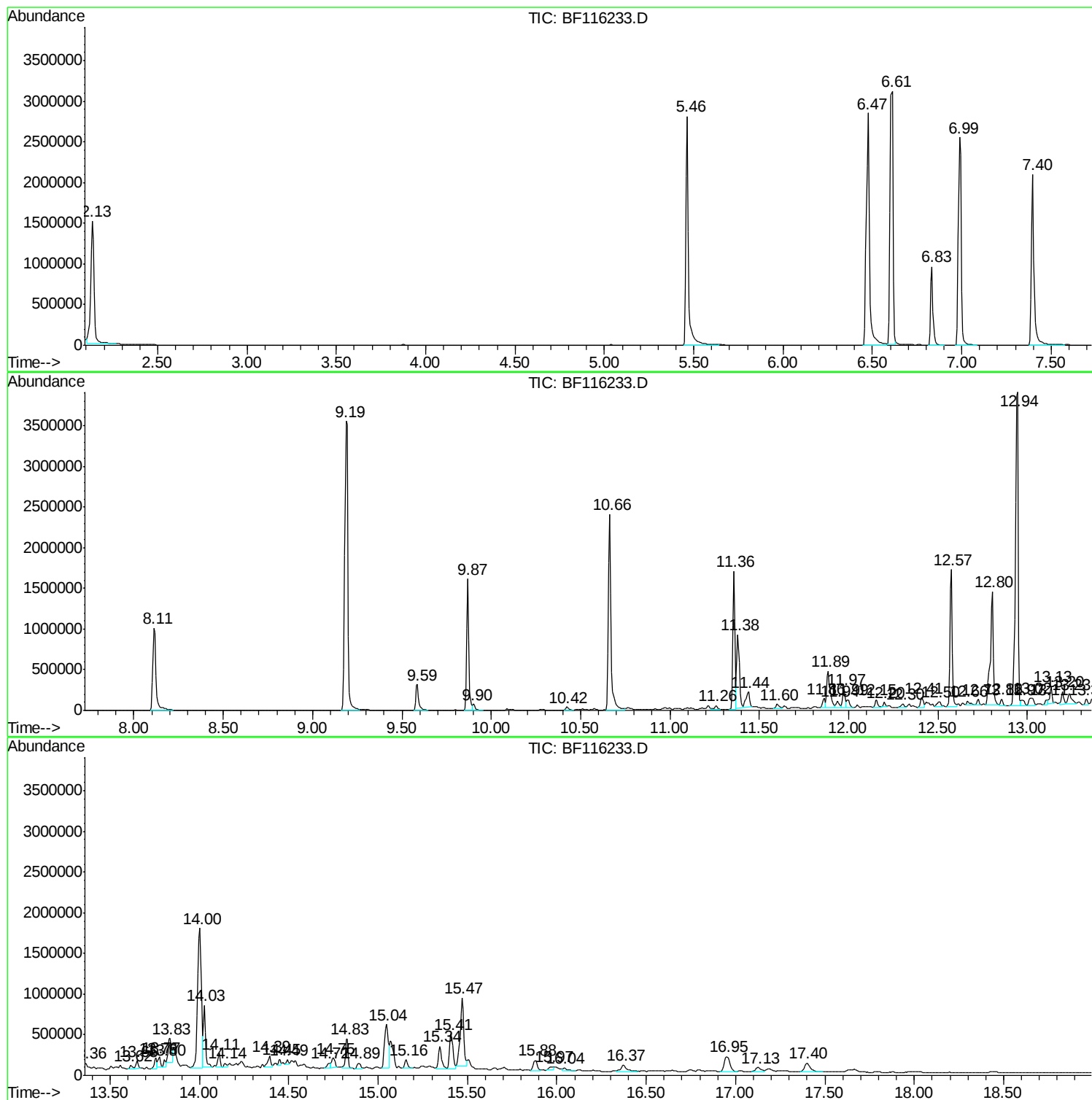
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ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
1S-I-(0-1)-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.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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Data File : BF116233.D
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Operator : HP/JU
Sample : K4227-21
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
1S-I-(0-1)-COMP

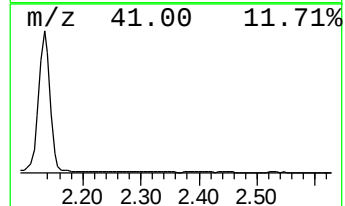
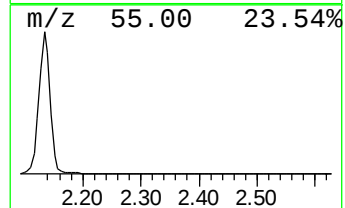
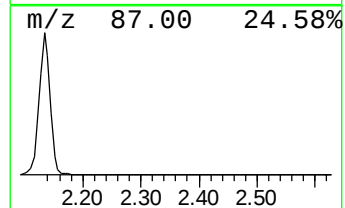
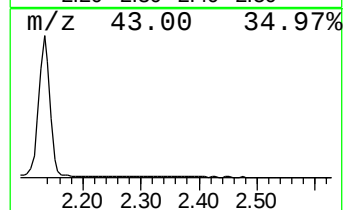
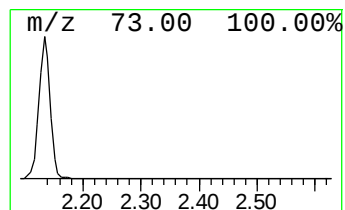
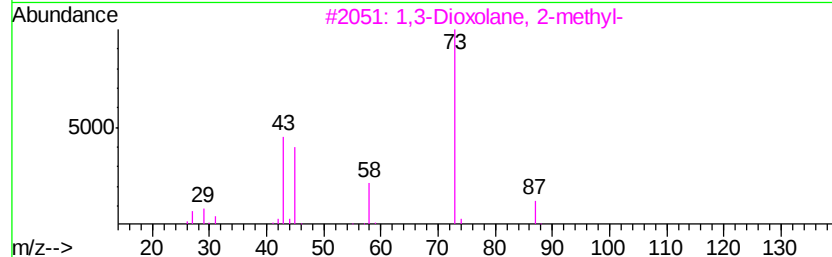
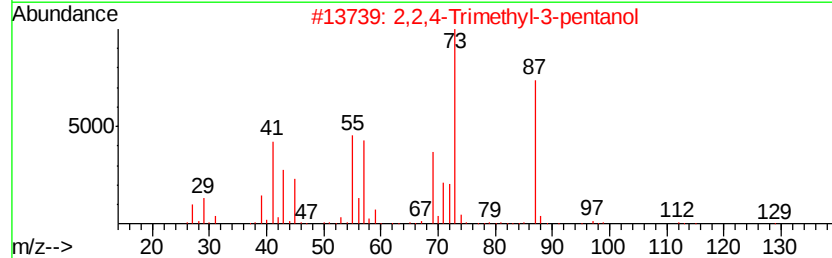
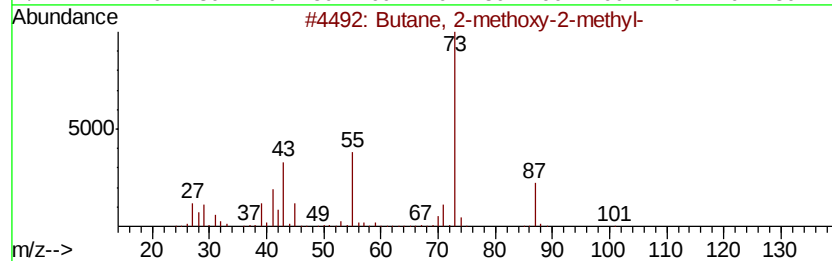
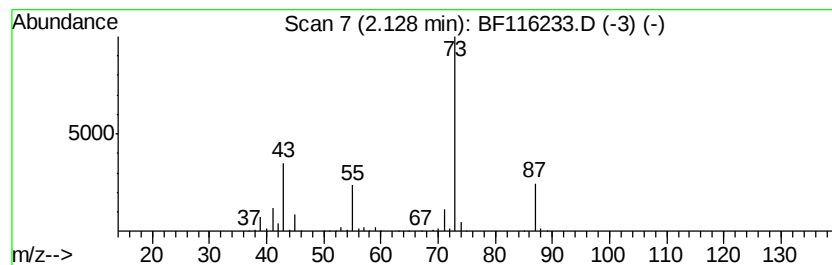
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
2.13	48.32 ng	2117240	1,4-Dichlorobenzene-d4	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	17



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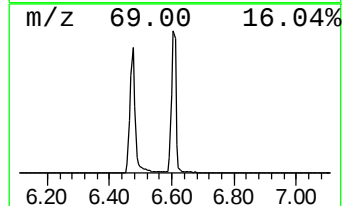
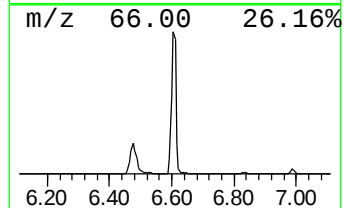
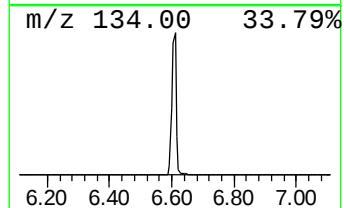
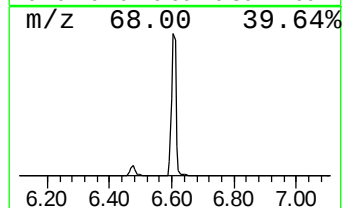
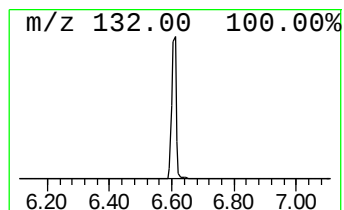
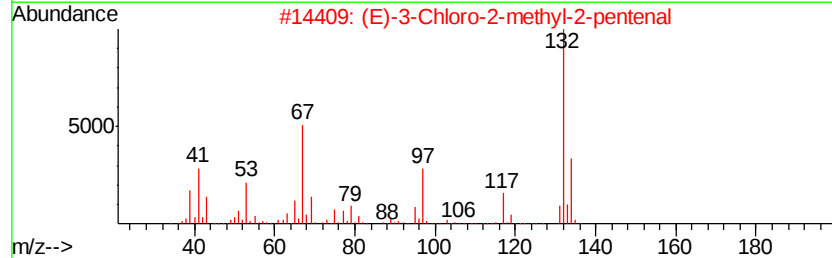
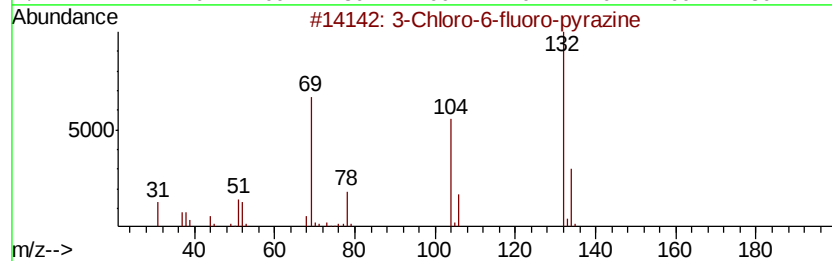
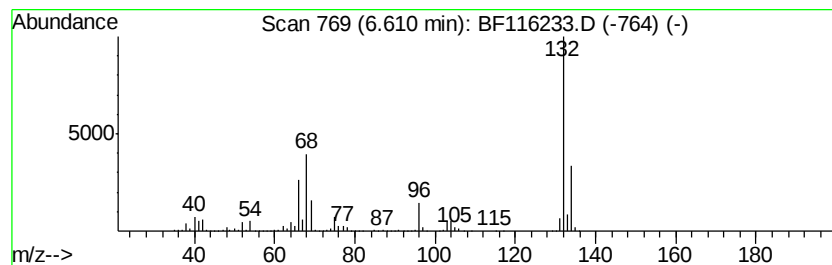
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.61	72.24 ng	3165350	1,4-Dichlorobenzene-d4	6.83

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		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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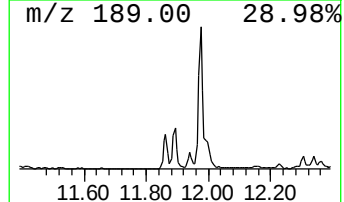
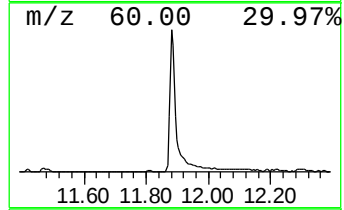
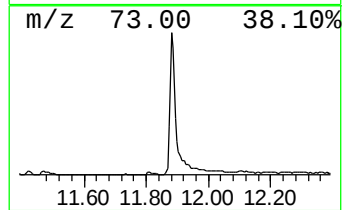
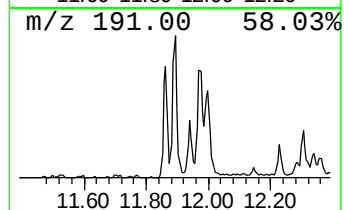
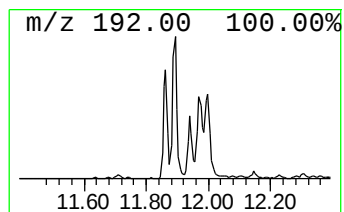
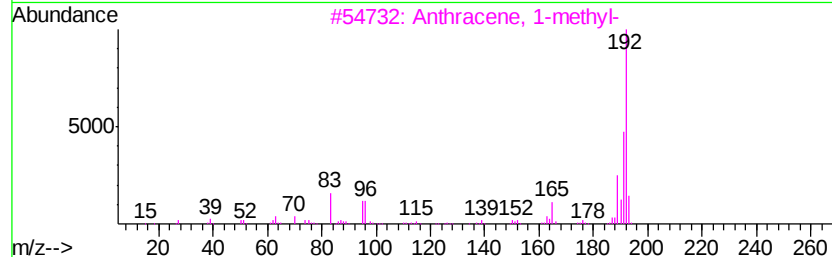
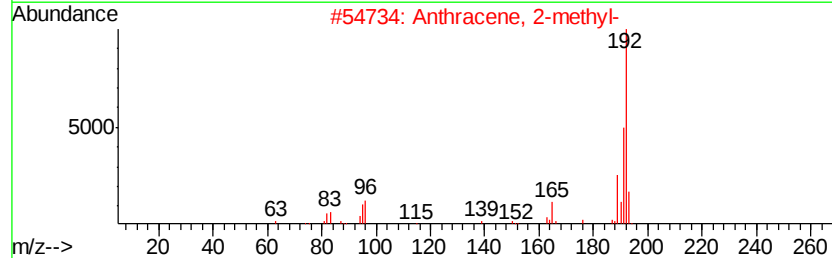
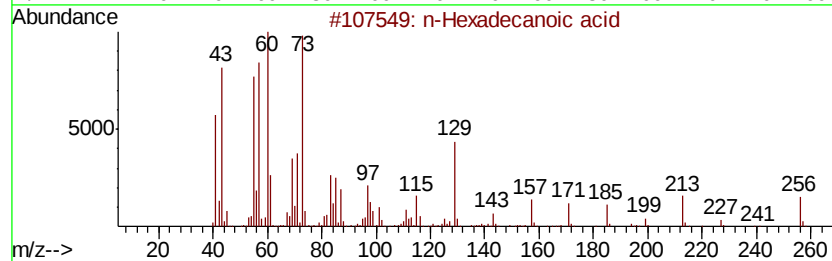
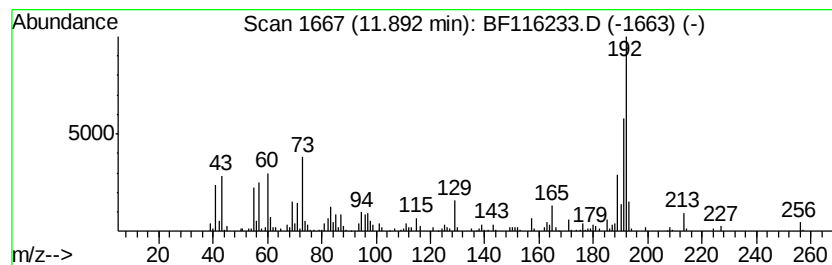
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.89	7.55 ng	583697	Phenanthrene-d10	11.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
3		Anthracene, 1-methyl-	192	C15H12	000610-48-0	91
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
5		Tridecanoic acid	214	C13H26O2	000638-53-9	89



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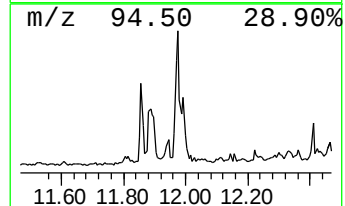
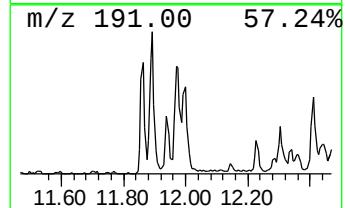
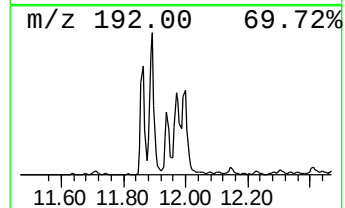
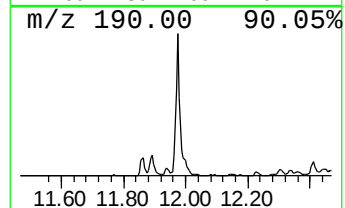
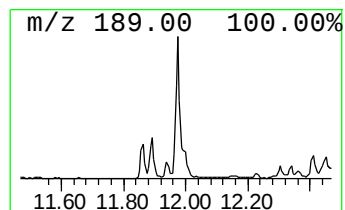
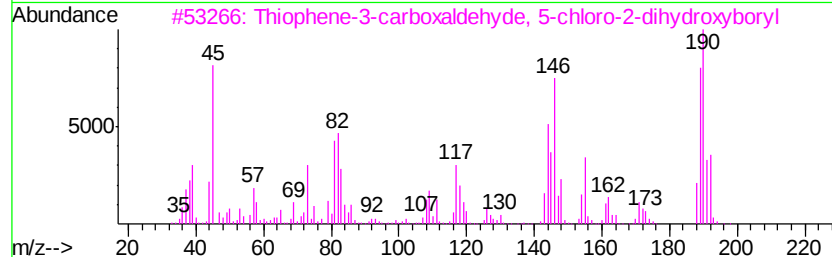
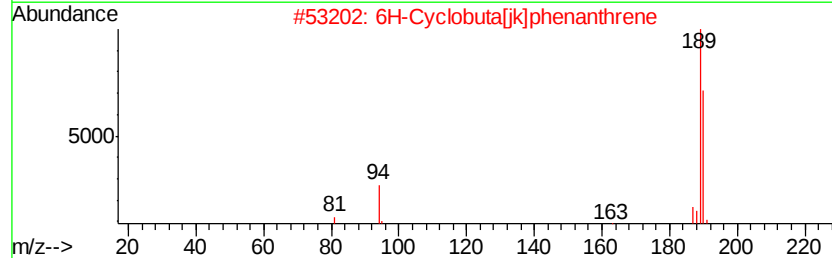
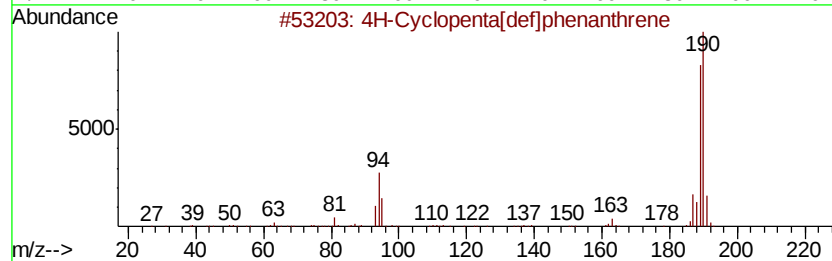
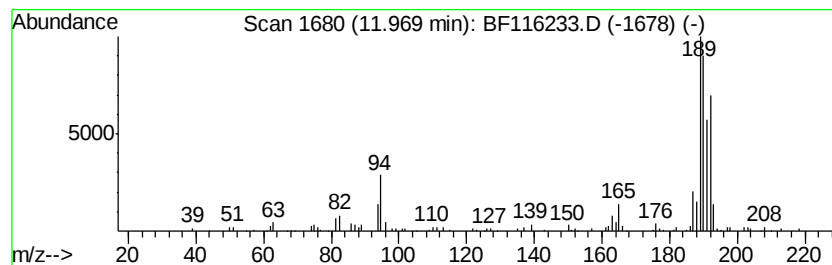
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BNA_F
ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.97	3.19 ng	246387	Phenanthrene-d10	11.36	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	81
2	6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	59
3	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	52
4	Methyl diselenide	190	C2H6Se2	007101-31-7	43
5	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081319\
Data File : BF116233.D
Acq On : 13 Aug 2019 15:54
Operator : HP/JU
Sample : K4227-21
Misc :
ALS Vial : 15 Sample Multiplier: 1

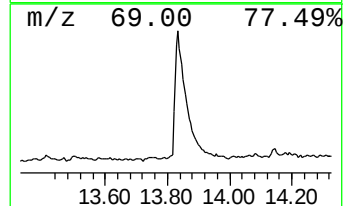
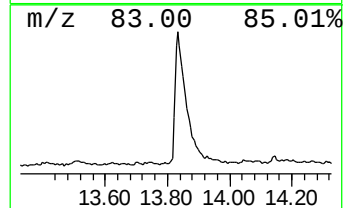
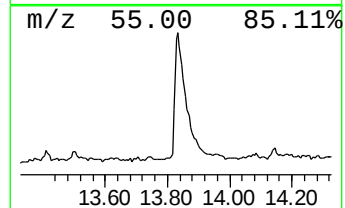
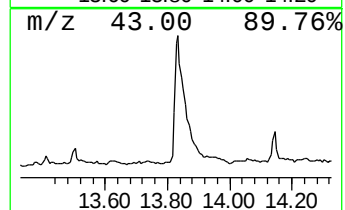
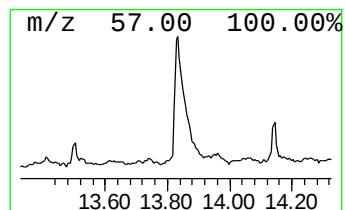
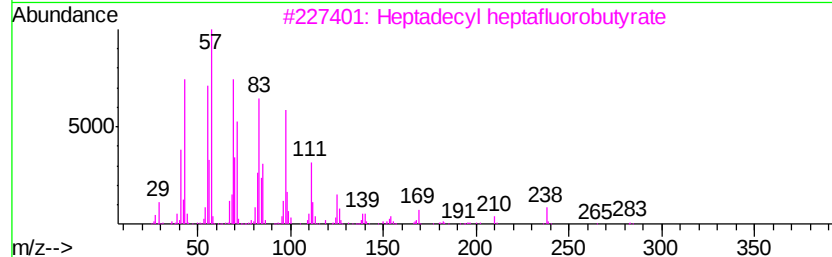
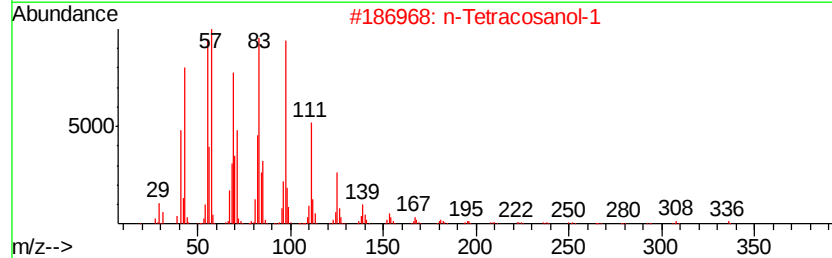
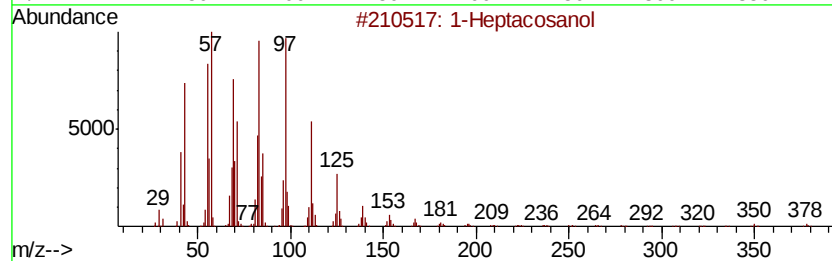
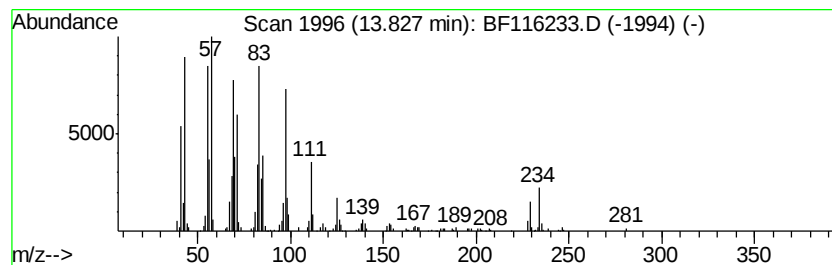
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 1-Heptacosanol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.83	3.67 ng	437735	Chrysene-d12	14.00	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heptacosanol	396	C27H56O	002004-39-9	93
2	n-Tetracosanol-1	354	C24H50O	000506-51-4	92
3	Heptadecyl heptafluorobutvrate	452	C21H35F7O2	959085-66-6	91
4	Behenic alcohol	326	C22H46O	000661-19-8	90
5	n-Heptadecanol-1	256	C17H36O	001454-85-9	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081319\
Data File : BF116233.D
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ClientSampleId :
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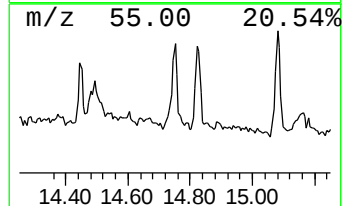
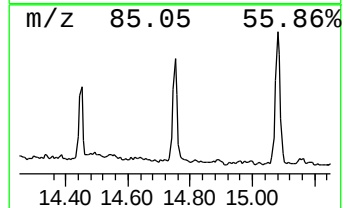
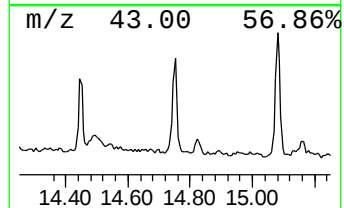
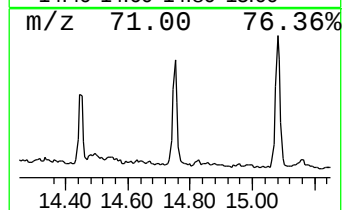
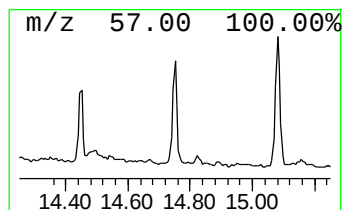
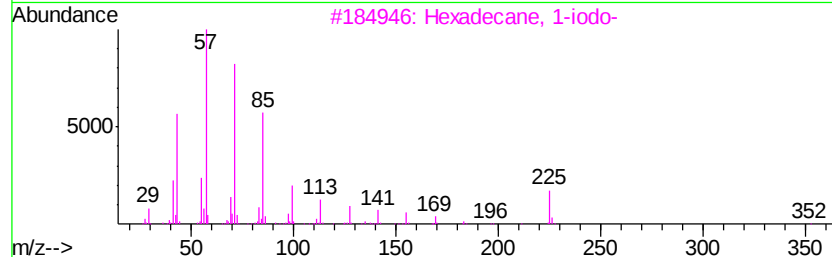
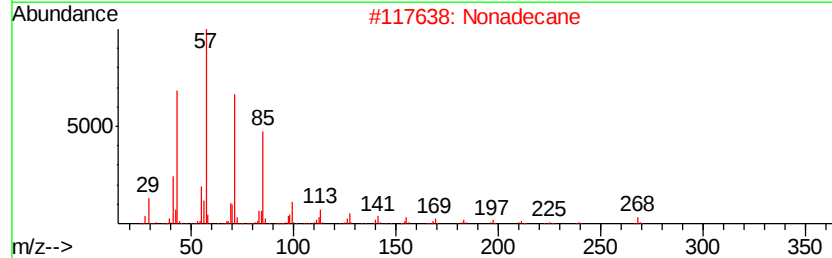
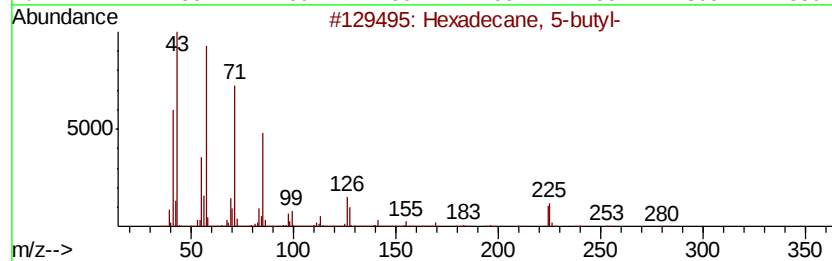
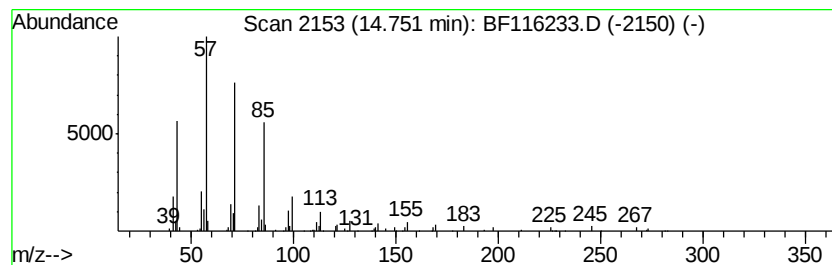
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexadecane, 5-butyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.75	2.10 ng	128021	Perylene-d12	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane, 5-butyl-	282	C20H42	006912-07-8	90
2		Nonadecane	268	C19H40	000629-92-5	86
3		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	86
4		Tetradecane, 4-ethyl-	226	C16H34	055045-14-2	86
5		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081319\
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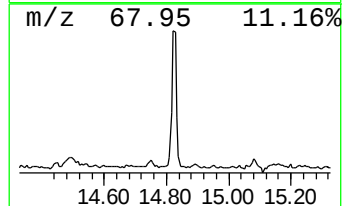
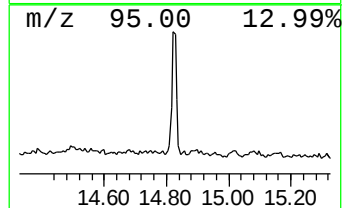
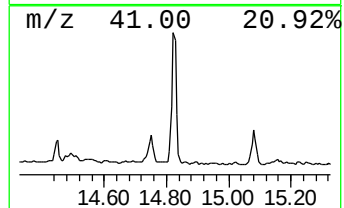
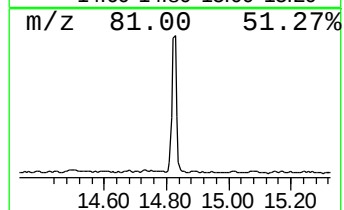
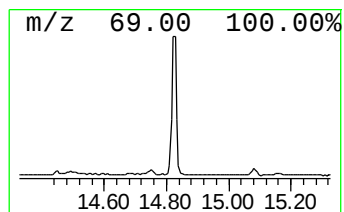
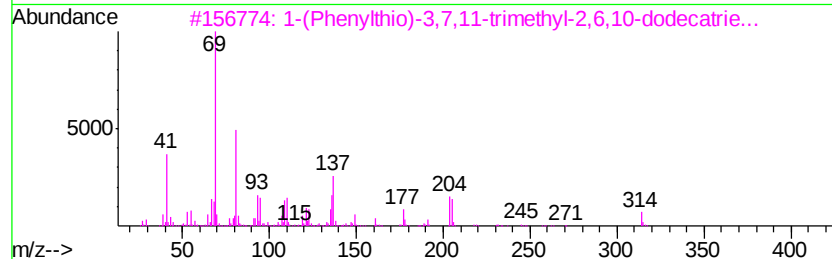
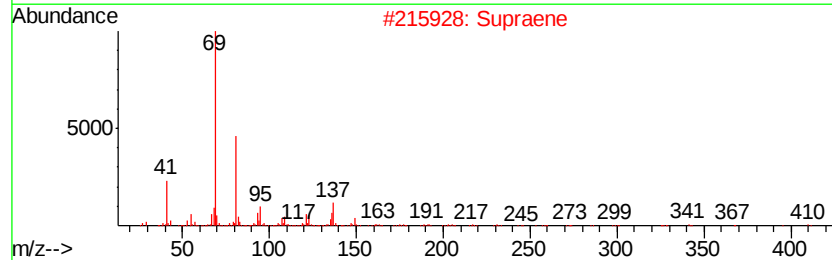
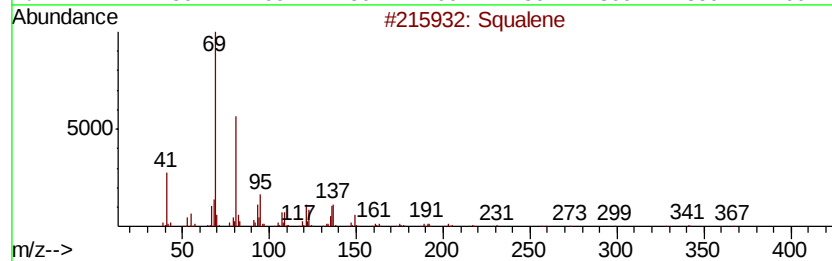
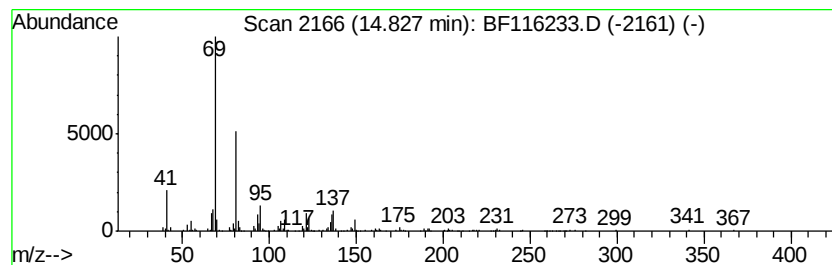
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Squalene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.		
14.83	6.52 ng	396992	Perylene-d12		15.47		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Squalene	410	C30H50	000111-02-4	94
2			Supraene	410	C30H50	007683-64-9	91
3			1-(Phenylthio)-3,7,11-trimethyl-...	314	C21H30S	028413-57-2	78
4			1,5,9-Decatriene, 2,3,5,8-tetram...	192	C14H24	230646-72-7	72
5			2,3,5-Trimethyl-2,3,5-hexanetric...	203	C12H17N3	003974-79-6	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081319\
Data File : BF116233.D
Acq On : 13 Aug 2019 15:54
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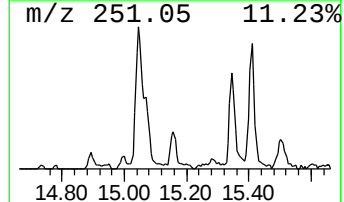
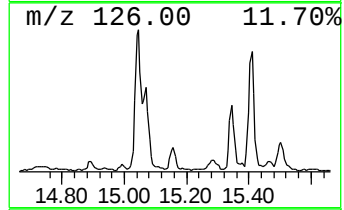
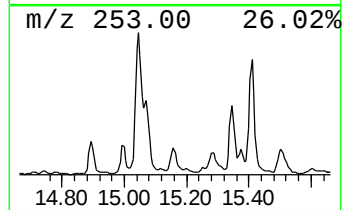
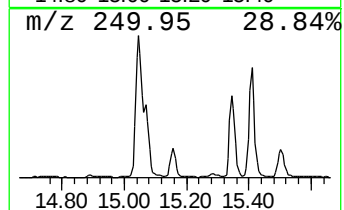
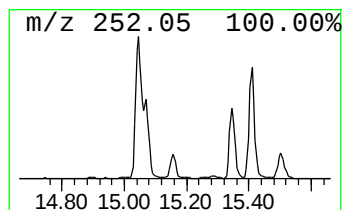
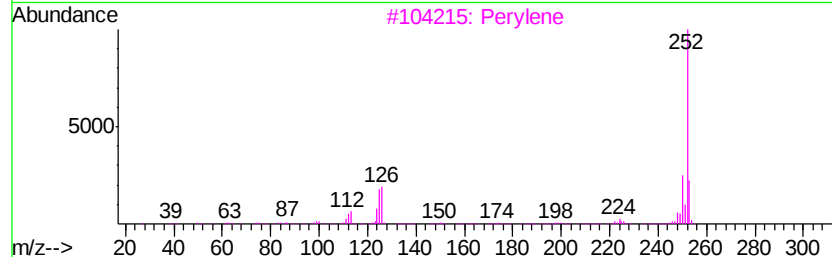
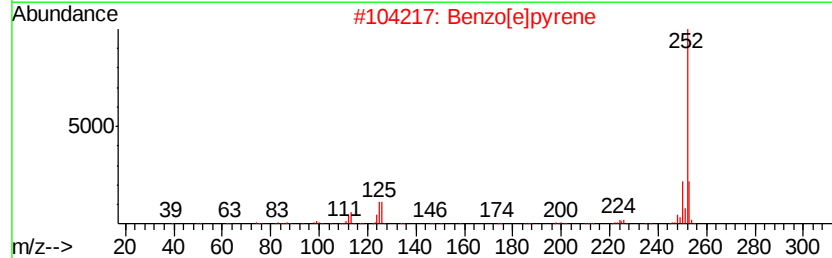
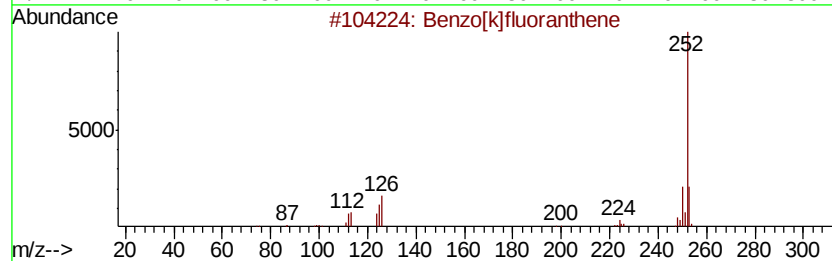
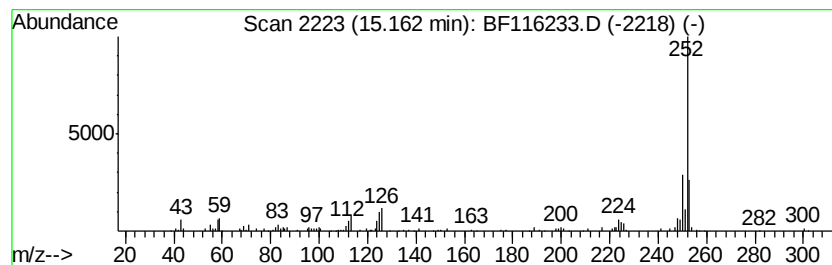
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzo[e]pyrene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.16	2.10 ng	128072	Perylene-d12	15.47

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzo[k]fluoranthene	252	C20H12	000207-08-9	98
2		Benzo[e]pyrene	252	C20H12	000192-97-2	96
3		Perylene	252	C20H12	000198-55-0	95
4		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	93
5		Benzo[a]pyrene	252	C20H12	000050-32-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081319\
Data File : BF116233.D
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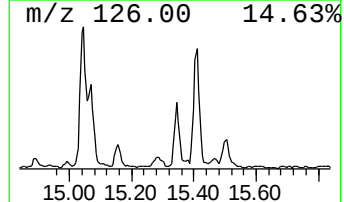
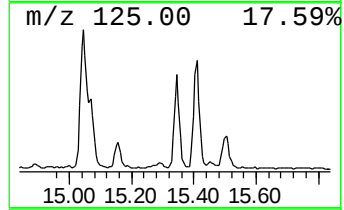
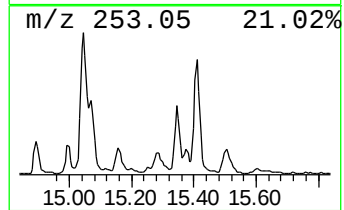
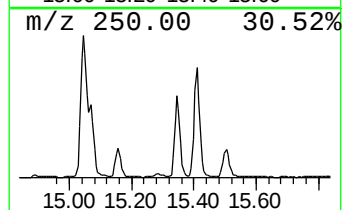
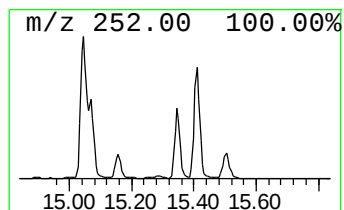
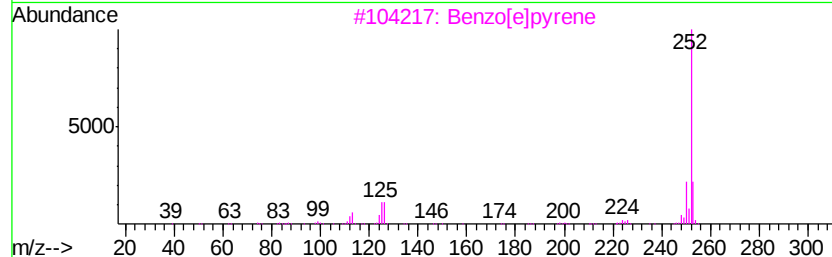
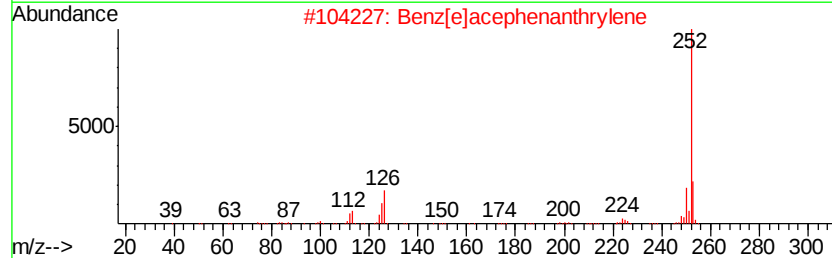
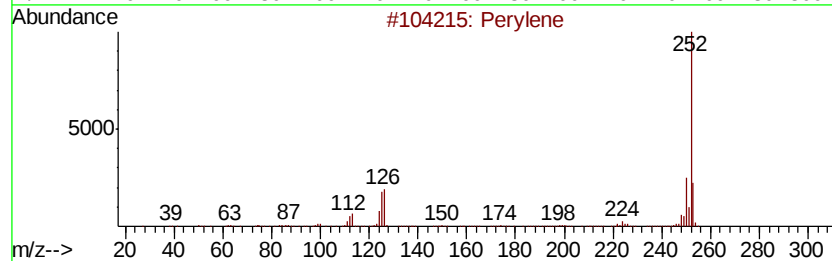
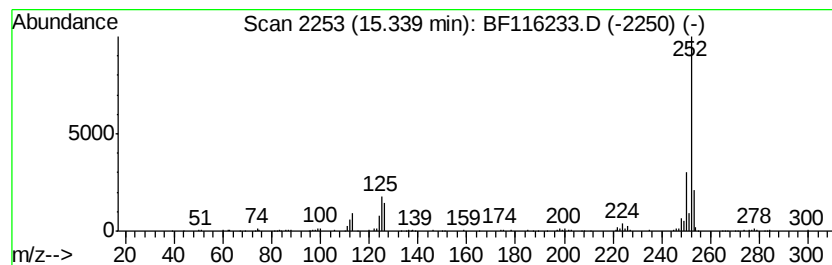
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Perylene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.34	5.83 ng	355026	Perylene-d12	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Perylene	252	C20H12	000198-55-0	98
2		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	98
3		Benzo[e]pyrene	252	C20H12	000192-97-2	98
4		Benzo[i]fluoranthene	252	C20H12	000205-82-3	97
5		Benzo[k]fluoranthene	252	C20H12	000207-08-9	97



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081319\
Data File : BF116233.D
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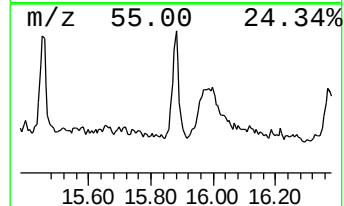
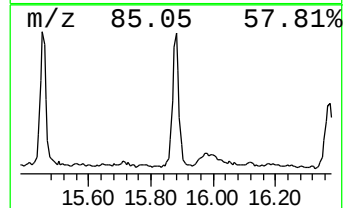
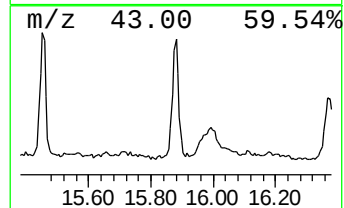
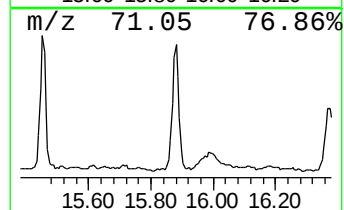
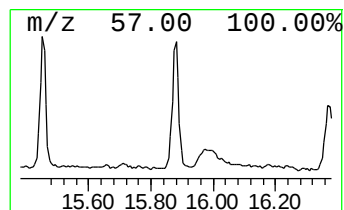
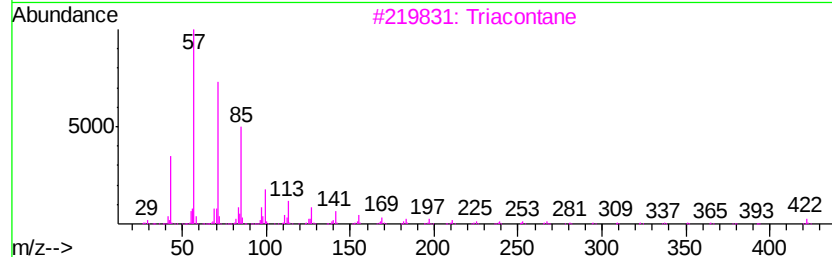
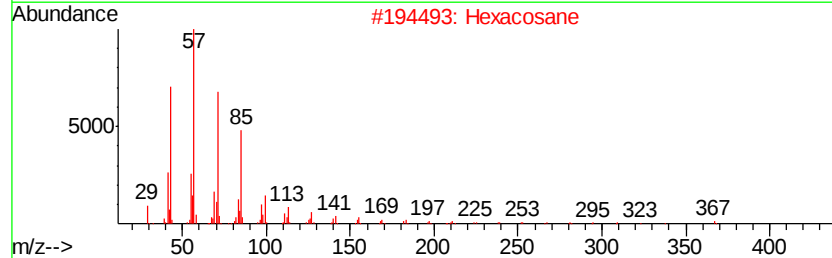
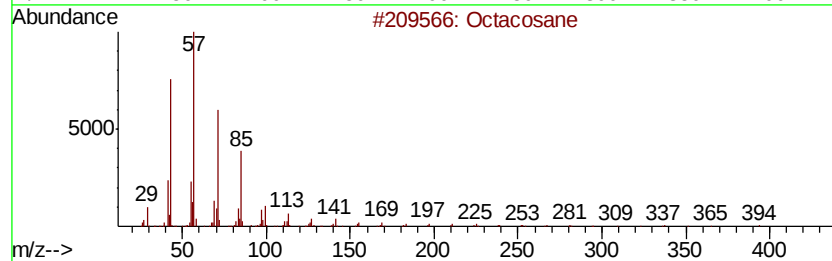
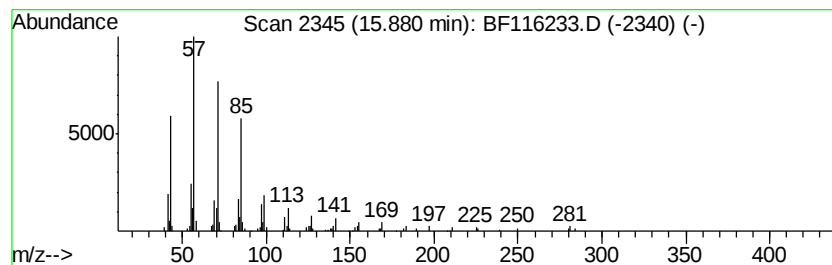
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octacosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.88	2.78 ng	169450	Perylene-d12	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Hexacosane	366	C26H54	000630-01-3	91
3		triacontane	422	C30H62	000638-68-6	91
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		Tetratetracontane	619	C44H90	007098-22-8	91



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ALS Vial : 15 Sample Multiplier: 1

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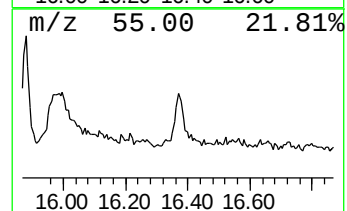
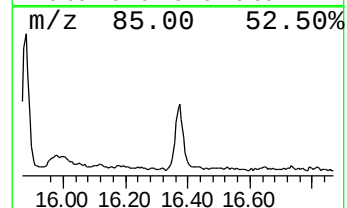
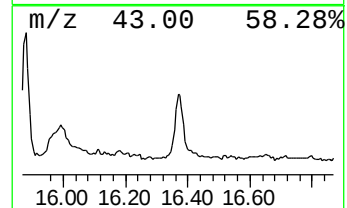
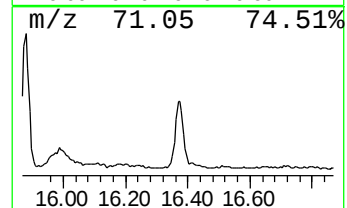
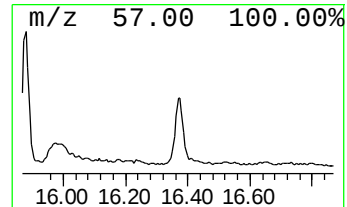
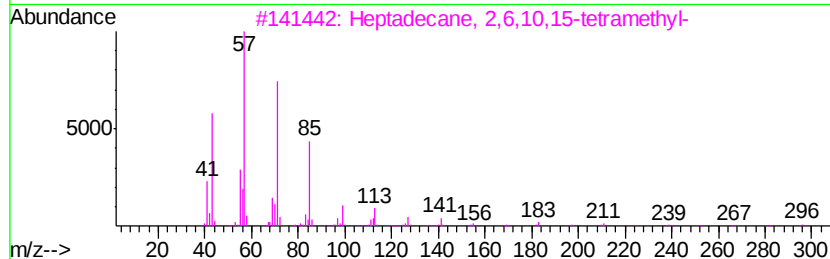
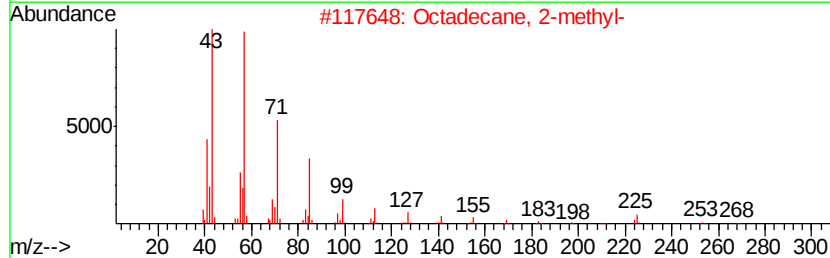
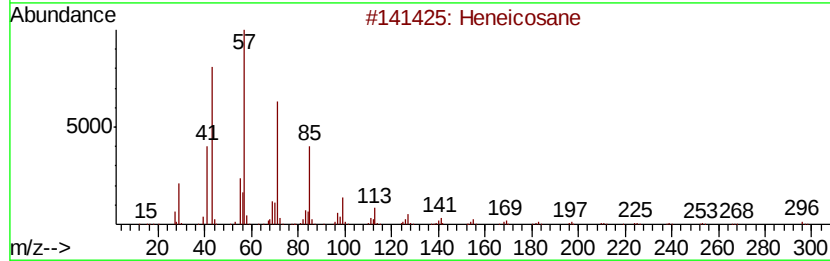
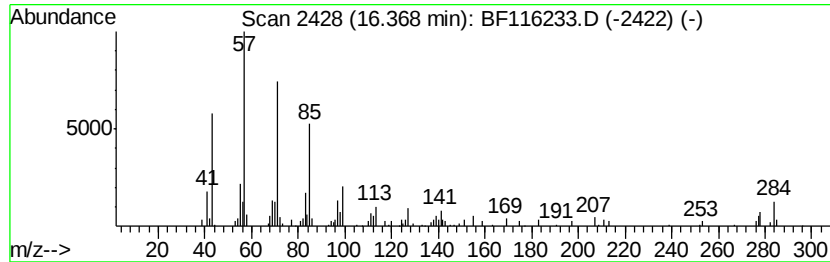
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heneicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.37	2.90 ng	176569	Perylene-d12	15.47

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	96
2			Octadecane, 2-methyl-	268	C19H40	001560-88-9	91
3			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	89
4			Triacontane	422	C30H62	000638-68-6	87
5			Hentriacontane	437	C31H64	000630-04-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF081319\
Data File : BF116233.D
Acq On : 13 Aug 2019 15:54
Operator : HP/JU
Sample : K4227-21
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
1S-I-(0-1)-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073019.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...	2.13	48.3	ng	2117240	1	6.83	876311	20.0
unknown6.61	6.61	72.2	ng	3165350	1	6.83	876311	20.0
n-Hexadecanoic acid	11.89	7.5	ng	583697	4	11.36	1546880	20.0
4H-Cyclopenta[def...	11.97	3.2	ng	246387	4	11.36	1546880	20.0
1-Heptacosanol	13.83	3.7	ng	437735	5	14.00	2385290	20.0
Hexadecane, 5-butyl-	14.75	2.1	ng	128021	6	15.47	1217560	20.0
Squalene	14.83	6.5	ng	396992	6	15.47	1217560	20.0
Benzo[e]pyrene	15.16	2.1	ng	128072	6	15.47	1217560	20.0
Perylene	15.34	5.8	ng	355026	6	15.47	1217560	20.0
Octacosane	15.88	2.8	ng	169450	6	15.47	1217560	20.0
Heneicosane	16.37	2.9	ng	176569	6	15.47	1217560	20.0