

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF052319\
 Data File : BF114545.D
 Acq On : 24 May 2019 1:06
 Operator : JU/SJ
 Sample : K2998-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EWING-SP

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-BF051019.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.439	567	570	590	rBV	2823474	3096358	34.04%	5.457%
2	6.457	739	743	761	rBV	3308111	3390549	37.28%	5.976%
3	6.586	761	765	772	rVV	3603278	3197751	35.16%	5.636%
4	6.810	800	803	812	rBV	1002156	882054	9.70%	1.555%
5	6.969	826	830	842	rBV	2589847	2266202	24.92%	3.994%
6	7.375	895	899	911	rBV	2313425	2040957	22.44%	3.597%
7	8.092	1017	1021	1030	rBV	1356494	1211765	13.32%	2.136%
8	9.169	1200	1204	1207	rBV	3758612	3237779	35.60%	5.706%
9	9.563	1267	1271	1277	rVB	452556	407298	4.48%	0.718%
10	9.845	1316	1319	1323	rVB	1440207	1324684	14.56%	2.335%
11	10.639	1450	1454	1465	rBV	2744597	2587911	28.45%	4.561%
12	11.333	1569	1572	1575	rBV	1687897	1626920	17.89%	2.867%
13	11.357	1575	1576	1580	rVB	396590	285266	3.14%	0.503%
14	11.863	1660	1662	1668	rVV3	306334	359979	3.96%	0.634%
15	11.910	1668	1670	1672	rVV2	169745	174851	1.92%	0.308%
16	11.933	1672	1674	1678	rVV2	463622	516548	5.68%	0.910%
17	11.980	1678	1682	1684	rVV	4297755	4170485	45.85%	7.350%
18	11.998	1684	1685	1688	rVB	184551	103593	1.14%	0.183%
19	12.033	1688	1691	1695	rBV6	57350	93510	1.03%	0.165%
20	12.080	1695	1699	1702	rBV	4787645	4182548	45.98%	7.371%
21	12.192	1714	1718	1724	rVB	2619668	2502828	27.52%	4.411%
22	12.551	1776	1779	1782	rBV	729600	607423	6.68%	1.071%
23	12.580	1782	1784	1788	rVB	323638	310667	3.42%	0.548%
24	12.780	1812	1818	1824	rBV	673092	731511	8.04%	1.289%
25	12.921	1837	1842	1845	rVB	3184729	3269560	35.95%	5.762%
26	13.074	1861	1868	1871	rBV	7257053	9095574	100.00%	16.030%
27	13.810	1990	1993	2002	rVB4	233505	421466	4.63%	0.743%
28	13.980	2018	2022	2025	rBV2	1442845	1514125	16.65%	2.669%
29	14.004	2025	2026	2029	rVB	267156	170895	1.88%	0.301%
30	14.733	2147	2150	2153	rVB	140162	139125	1.53%	0.245%
31	15.021	2196	2199	2202	rBV	145008	152256	1.67%	0.268%
32	15.321	2247	2250	2255	rVB	152473	188412	2.07%	0.332%
33	15.380	2257	2260	2264	rVB	203043	248060	2.73%	0.437%
34	15.445	2265	2271	2276	rBV	966502	1416312	15.57%	2.496%

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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-BF051019.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	15.856	2337	2341	2346	rVB	220539	304215	3.34%	0.536%
36	16.345	2420	2424	2430	rBV2	143335	274949	3.02%	0.485%
37	16.915	2518	2521	2527	rVB2	130909	235288	2.59%	0.415%

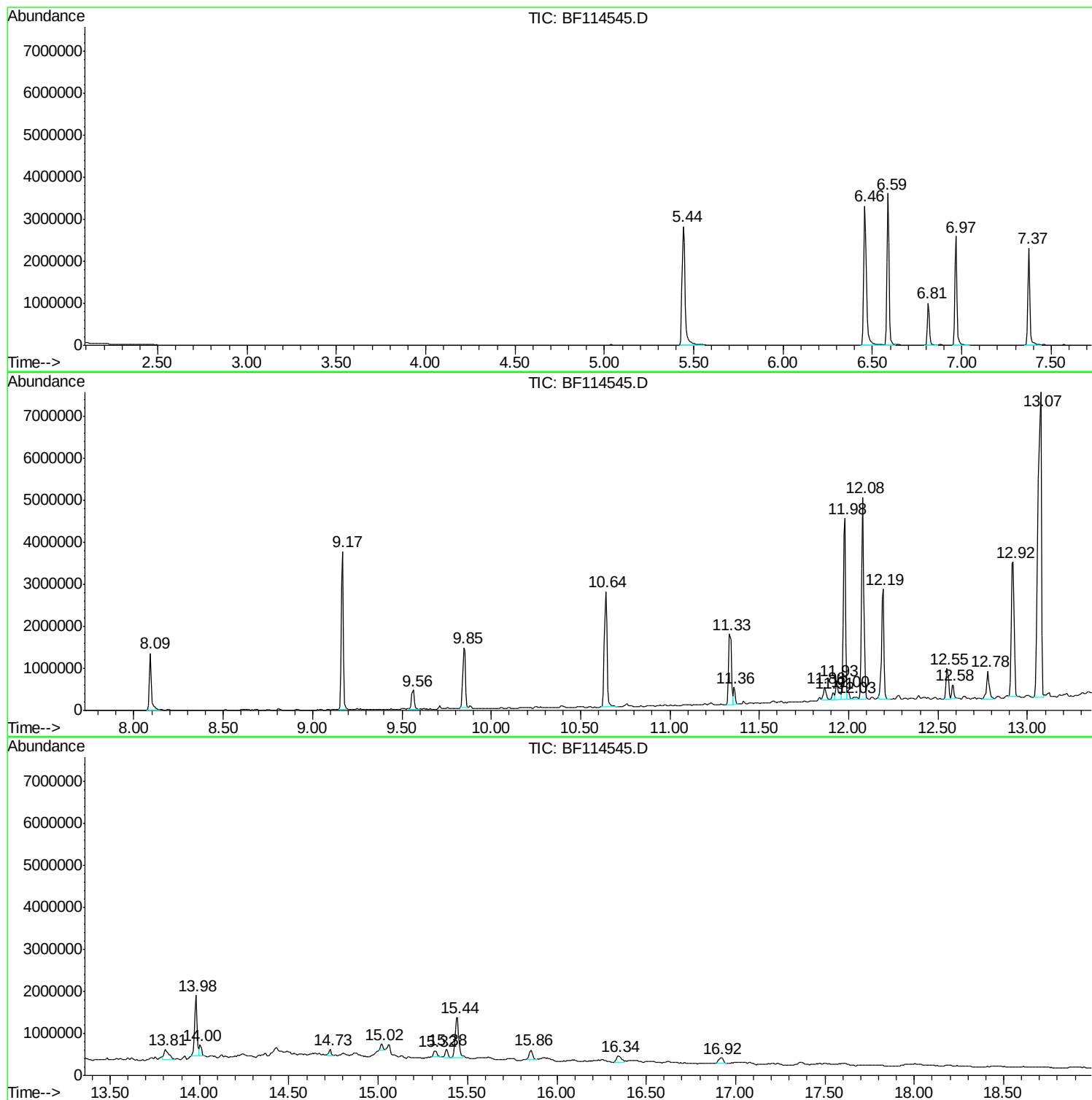
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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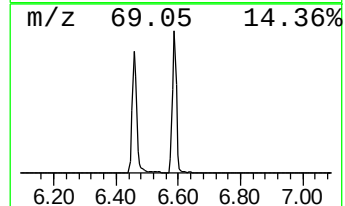
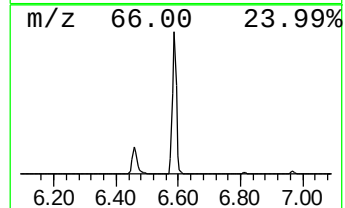
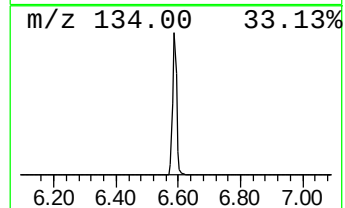
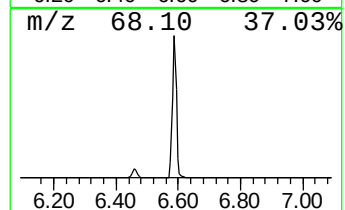
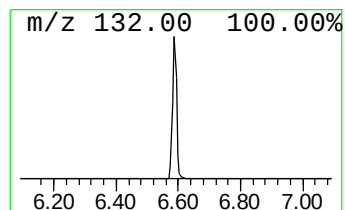
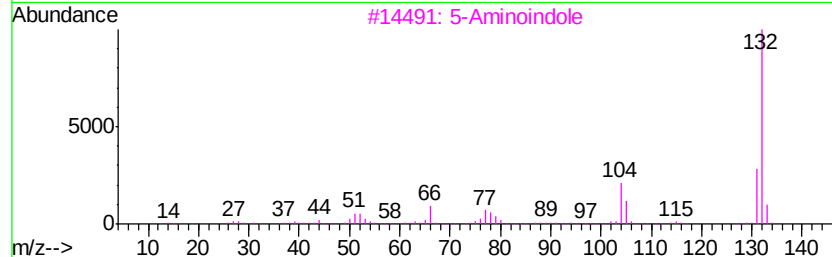
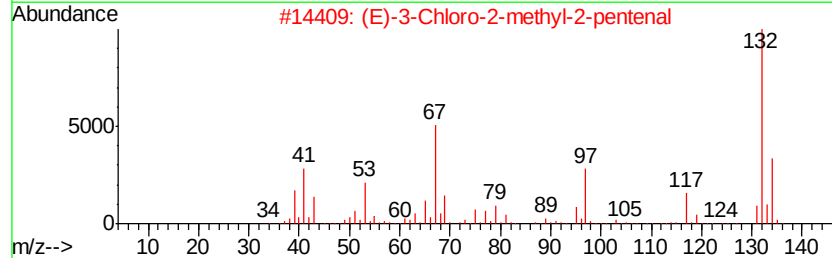
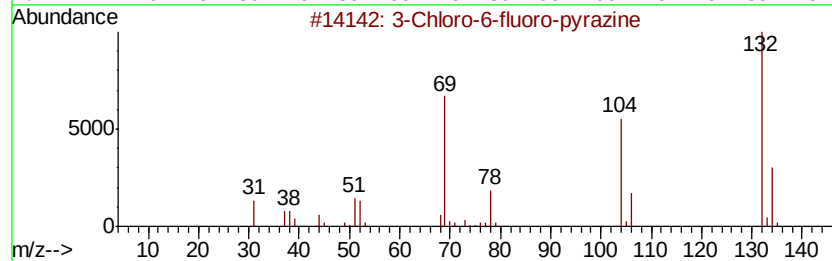
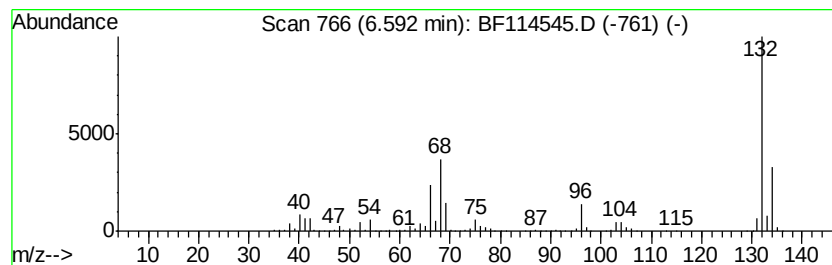
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.59 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.59	72.51 ng	3197750	1,4-Dichlorobenzene-d4	6.81

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	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



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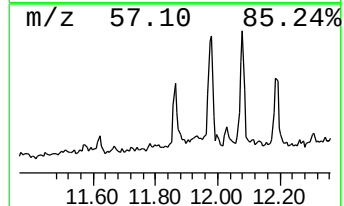
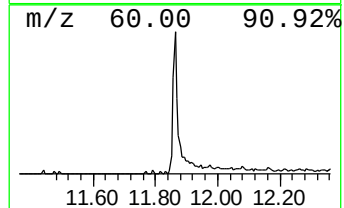
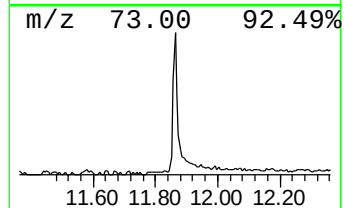
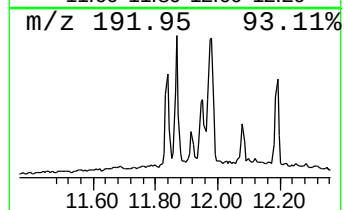
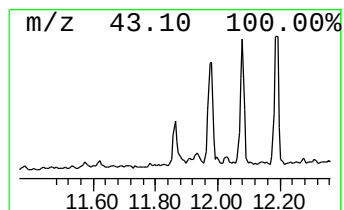
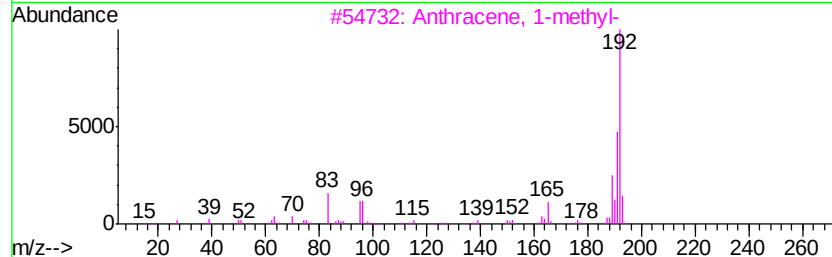
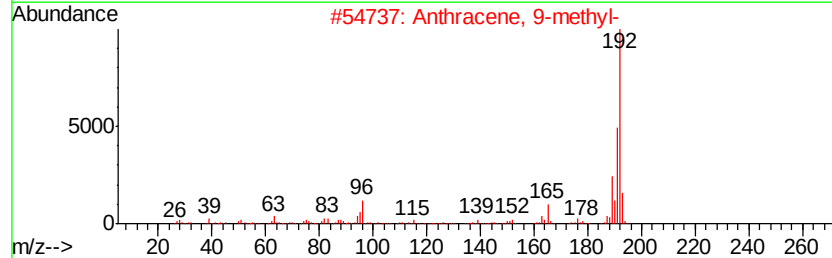
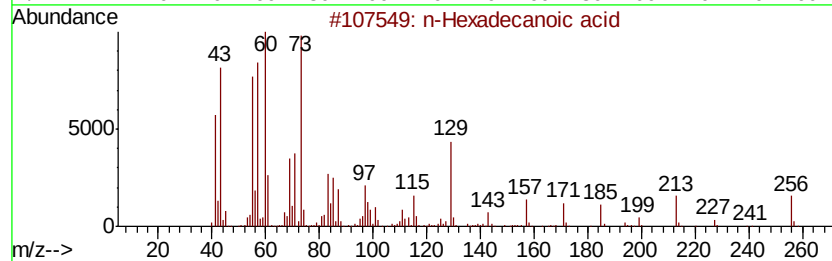
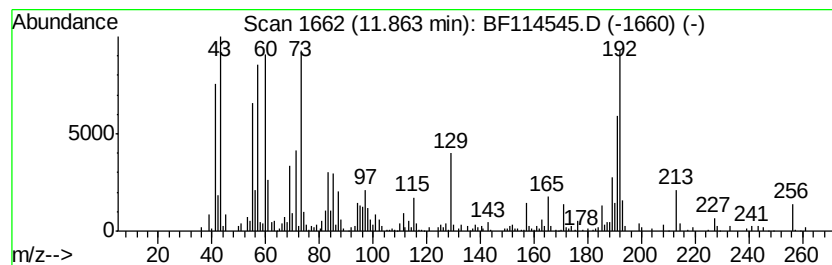
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 3 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.86	4.43 ng	359979	Phenanthrene-d10		11.33
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Anthracene, 9-methyl-	192	C15H12	000779-02-2	68
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	68
4	1H-Cyclopropa[1,1]phenanthrene,1a,...	192	C15H12	000949-41-7	68
5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	68



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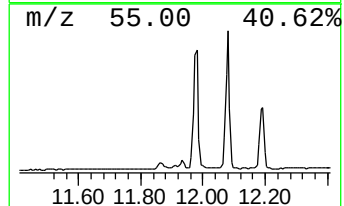
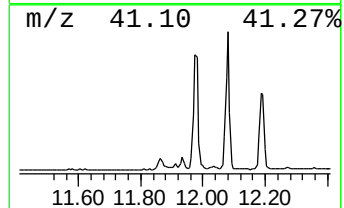
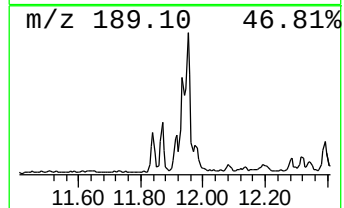
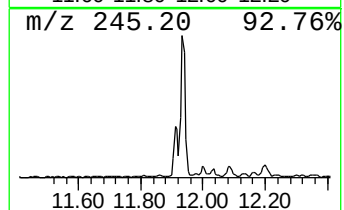
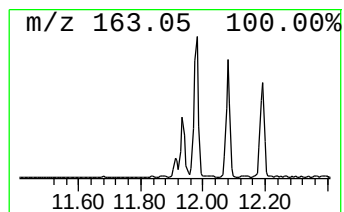
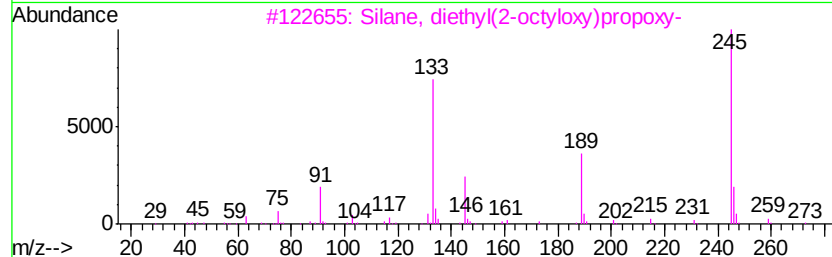
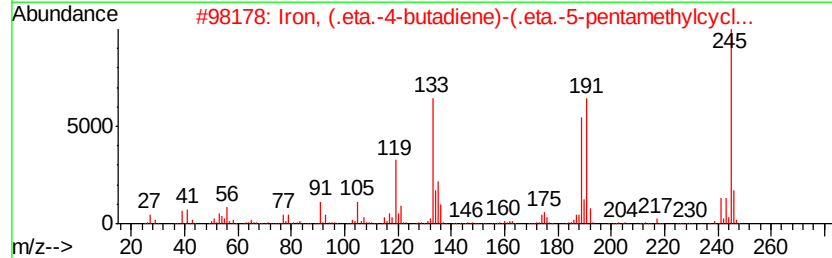
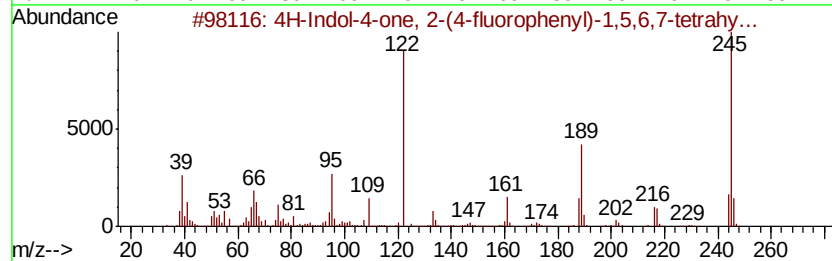
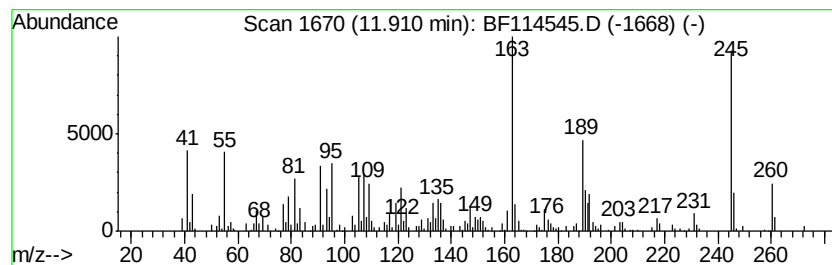
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 unknown11.91 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.91	2.15 ng	174851	Phenanthrene-d10	11.33	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Indol-4-one, 2-(4-fluoropheny...	245	C14H12FN02	1000337-74-5	45
2	Iron, (.eta.-4-butadiene)-(.eta....	245	C14H21Fe	1000161-69-5	27
3	Silane, diethyl(2-octvloxy)propoxv-	274	C15H3402Si	1000363-00-9	22
4	Silane, dimethylnonyloxypropyloxy-	260	C14H3202Si	1000356-04-5	14
5	Alpha-(4-cinnolinyl)-alpha-pheny...	245	C16H11N3	018592-05-7	14



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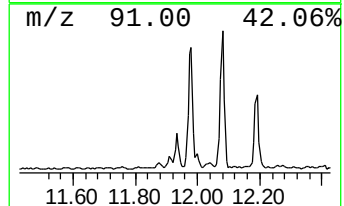
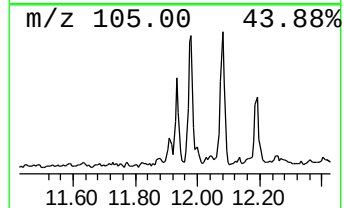
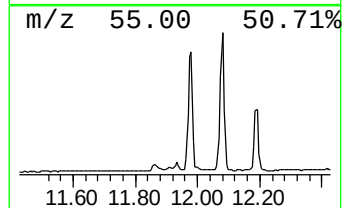
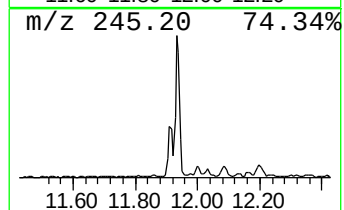
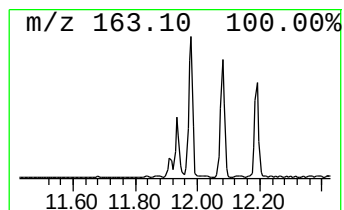
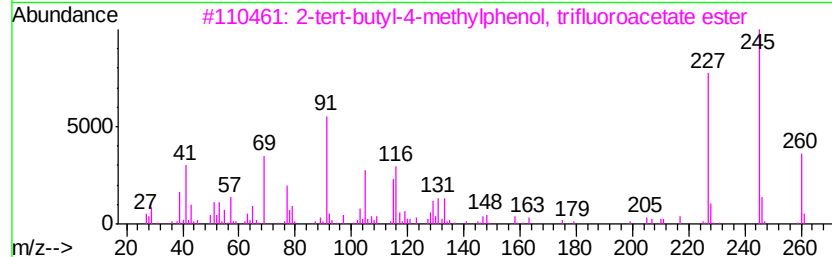
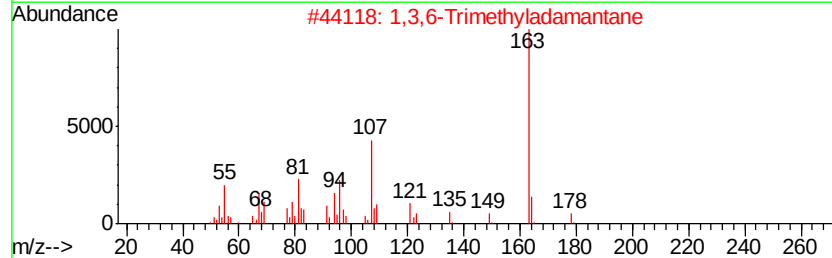
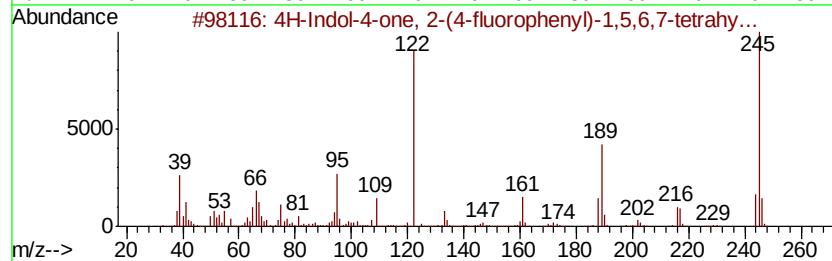
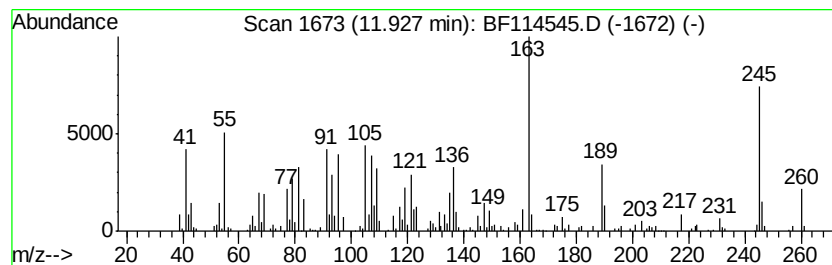
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 unknown11.93 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.93	6.35 ng	516548	Phenanthrene-d10	11.33	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Indol-4-one, 2-(4-fluorophenv...	245	C14H12FN02	1000337-74-5	38
2	1,3,6-Trimethyladamantane	178	C13H22	024139-37-5	35
3	2-tert-butvl-4-methylphenol, tri...	260	C13H15F302	1000376-40-6	30
4	N-Acetyl-L-tyrosinamide	222	C11H14N203	001948-71-6	22
5	1,3,4-Trimethyladamantane	178	C13H22	1000214-98-3	22



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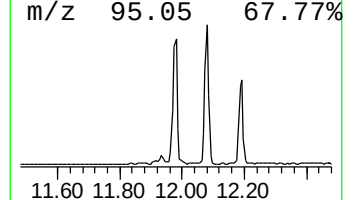
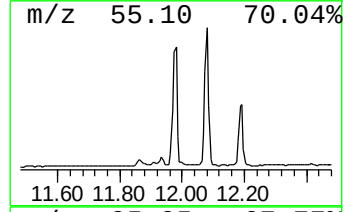
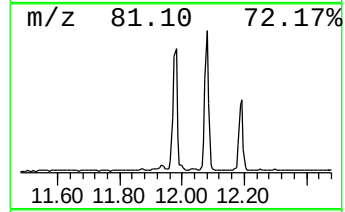
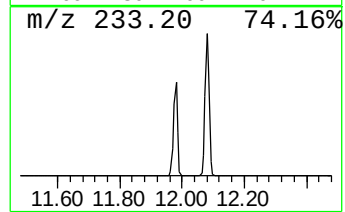
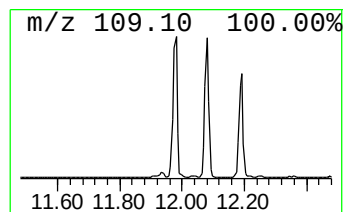
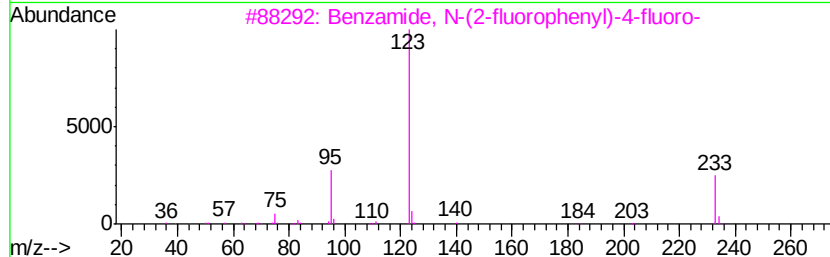
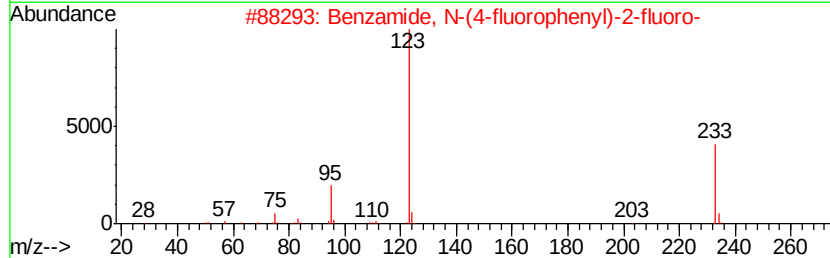
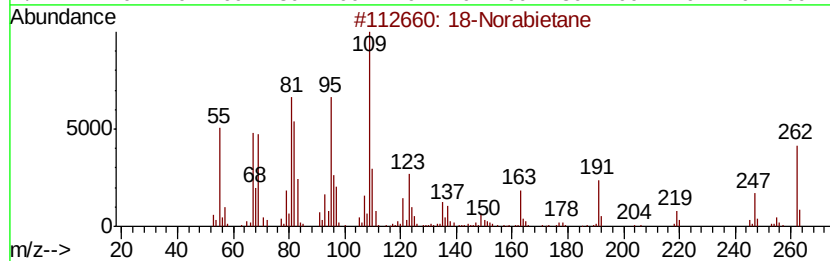
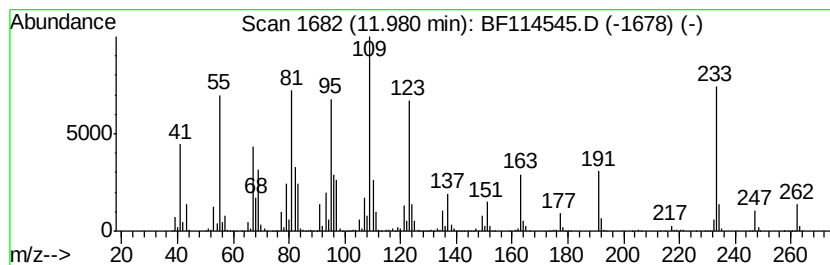
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 18-Norabietane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.98	51.27 ng	4170490	Phenanthrene-d10	11.33	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	18-Norabietane	262	C19H34	1000293-16-6	70
2	Benzamide, N-(4-fluorophenyl)-2-...	233	C13H9F2NO	1000307-07-9	43
3	Benzamide, N-(2-fluorophenyl)-4-...	233	C13H9F2NO	1000308-30-2	38
4	Benzamide, N-(4-fluorophenyl)-3-...	233	C13H9F2NO	1000307-16-3	38
5	Benzamide, N-(2-fluorophenyl)-2-...	233	C13H9F2NO	1000308-09-1	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052319\
Data File : BF114545.D
Acq On : 24 May 2019 1:06
Operator : JU/SJ
Sample : K2998-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EWING-SP

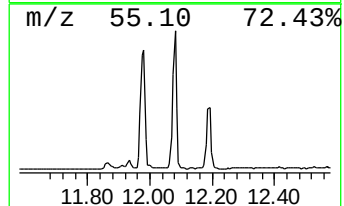
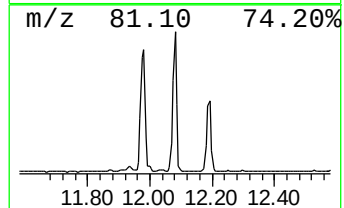
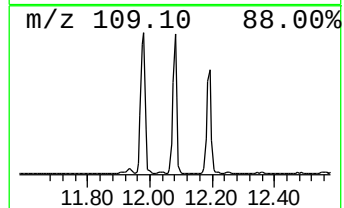
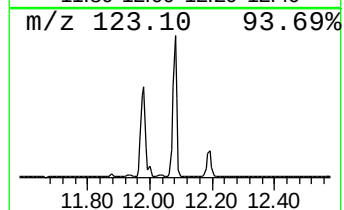
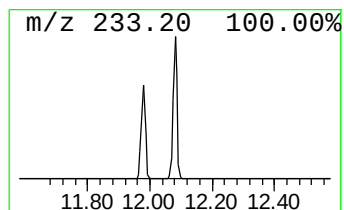
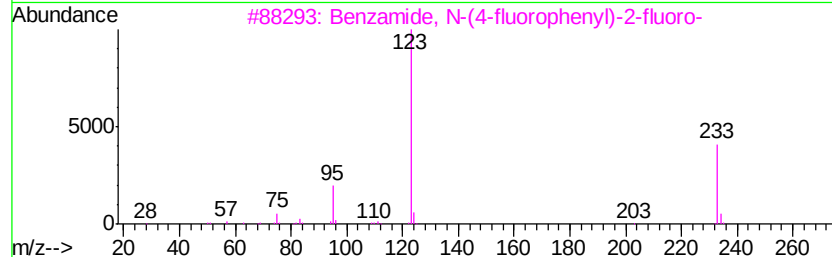
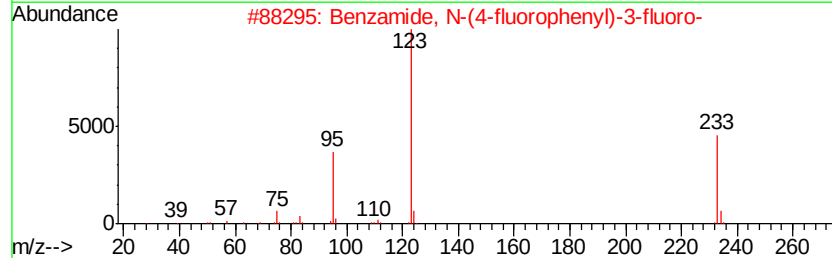
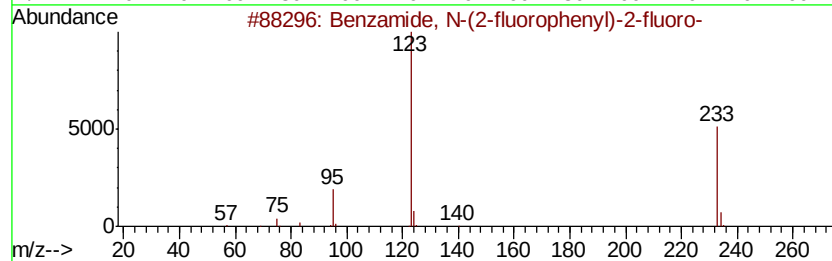
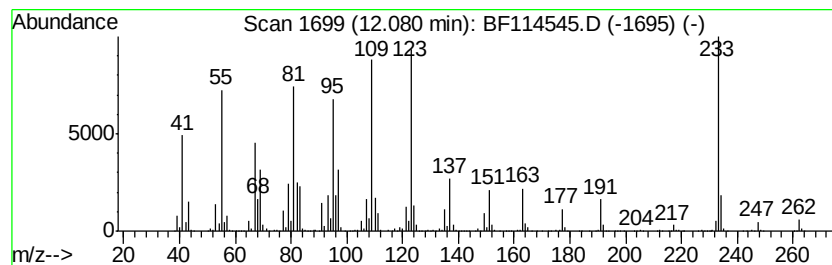
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown12.08 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.08	51.42 ng	4182550	Phenanthrene-d10	11.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzamide, N-(2-fluorophenyl)-2-...	233	C13H9F2NO	1000308-09-1	38
2		Benzamide, N-(4-fluorophenyl)-3-...	233	C13H9F2NO	1000307-16-3	38
3		Benzamide, N-(4-fluorophenyl)-2-...	233	C13H9F2NO	1000307-07-9	35
4		trans-Chrysanthemal	152	C10H16O	020104-05-6	30
5		Ethyl chrysanthemate	196	C12H20O2	000097-41-6	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052319\
Data File : BF114545.D
Acq On : 24 May 2019 1:06
Operator : JU/SJ
Sample : K2998-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

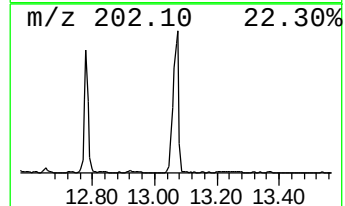
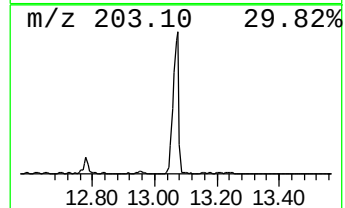
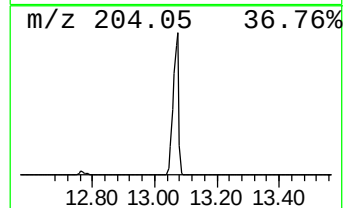
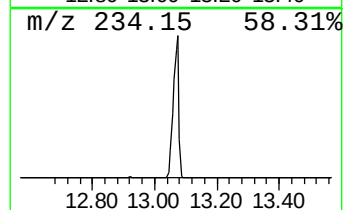
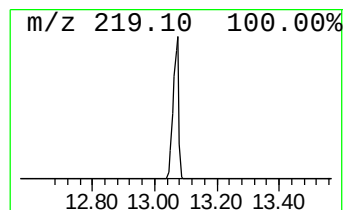
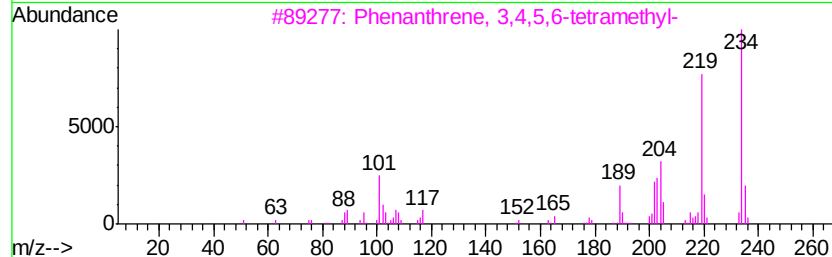
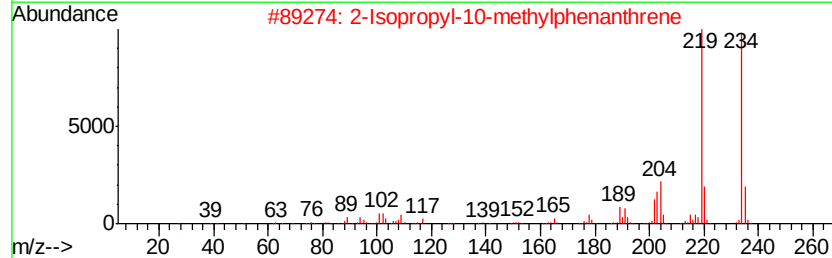
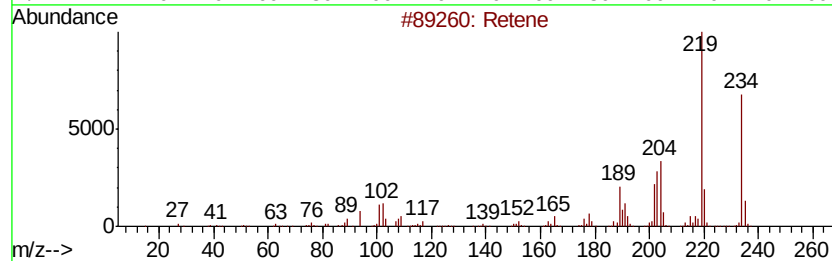
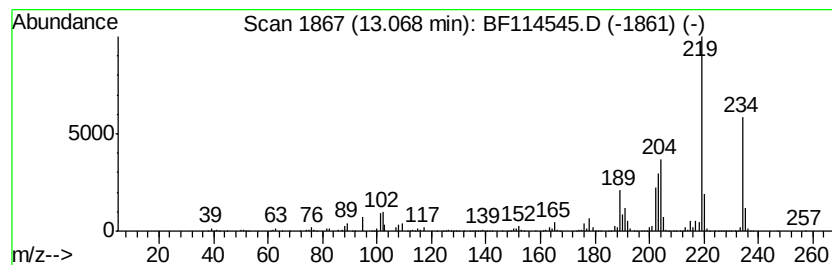
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Retene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.07	120.14 ng	9095570	Chrysene-d12	13.98	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Retene	234	C18H18	000483-65-8	99
2	2-Isopropyl-10-methylphenanthrene	234	C18H18	066552-97-4	95
3	Phenanthrene, 3,4,5,6-tetramethyl-	234	C18H18	007343-06-8	70
4	3,5-di-tert-Butyl-4-hydroxybenza...	234	C15H22O2	001620-98-0	50
5	2,3,5,6-Detetrahydrocyclohexanon...	234	C15H22O2	101100-38-3	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052319\
Data File : BF114545.D
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Sample : K2998-01
Misc :
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ClientSampleId :
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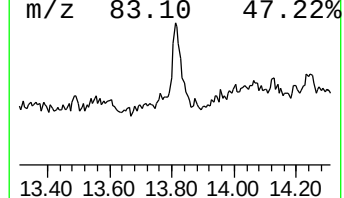
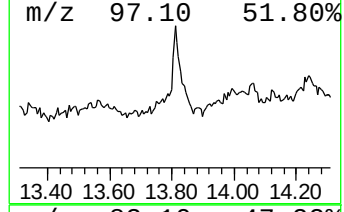
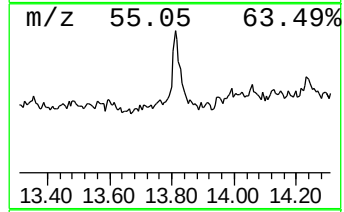
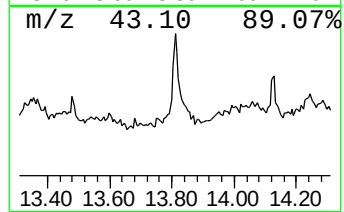
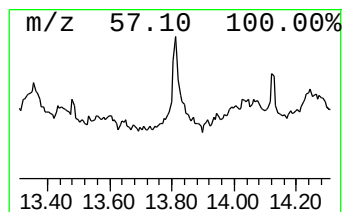
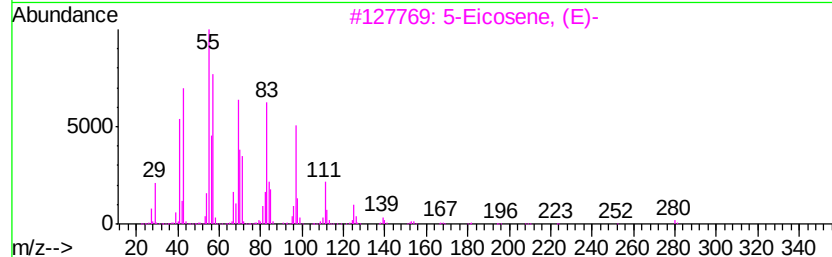
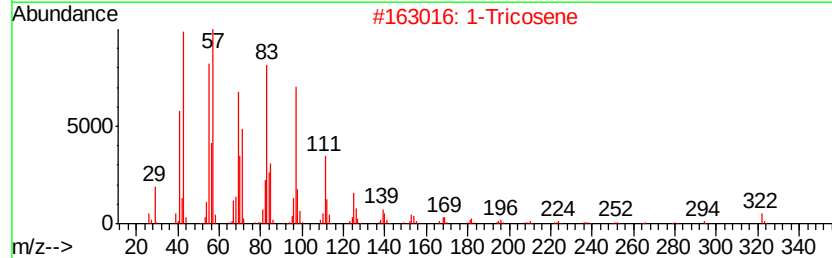
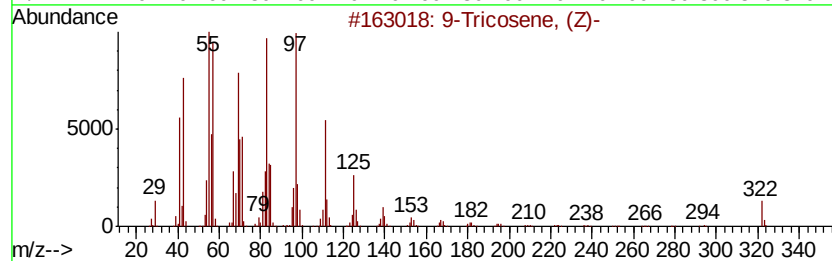
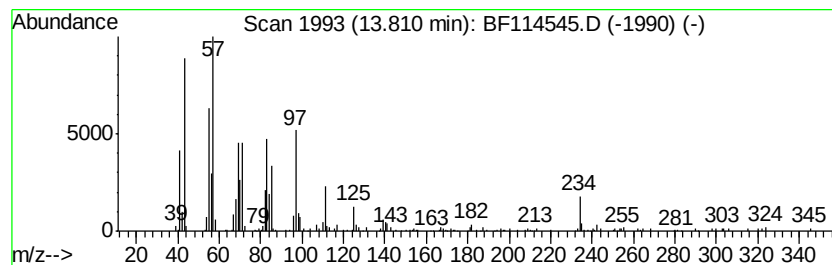
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 9-Tricosene, (Z)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.81	5.57 ng	421466	Chrysene-d12	13.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Tricosene, (Z)-	322	C23H46	027519-02-4	90
2		1-Tricosene	322	C23H46	018835-32-0	90
3		5-Eicosene, (E)-	280	C20H40	074685-30-6	90
4		1-Hexadecanesulfonyl chloride	324	C16H33ClO2S	038775-38-1	87
5		Eicosyl heptafluorobutyrate	494	C24H41F7O2	1000351-83-5	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052319\
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Sample : K2998-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

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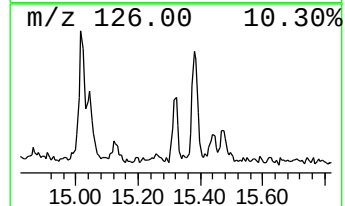
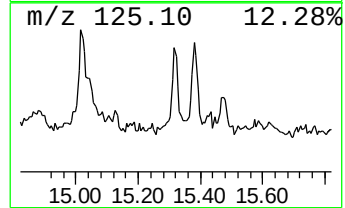
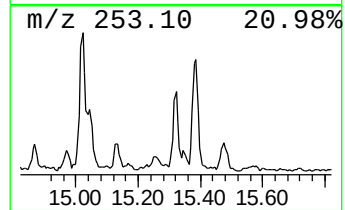
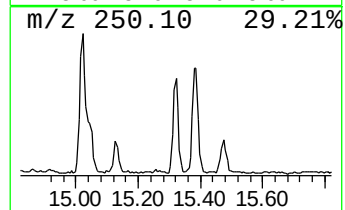
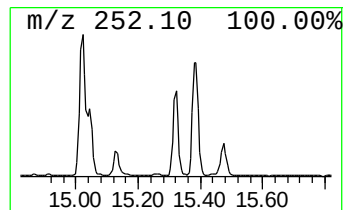
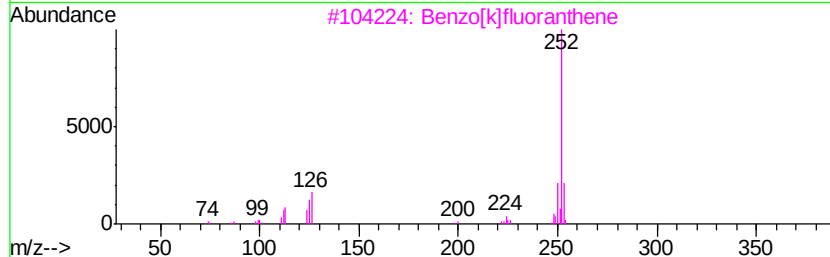
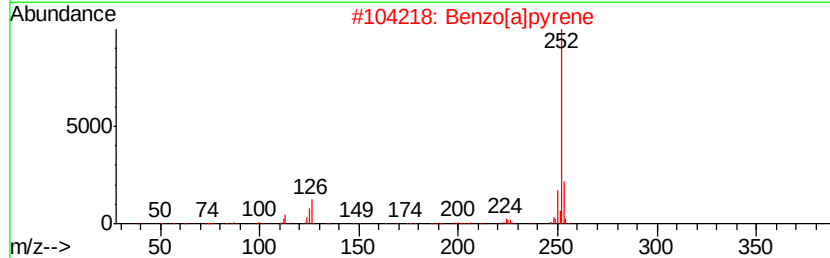
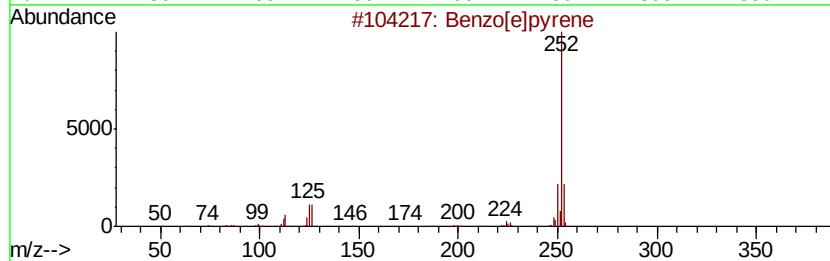
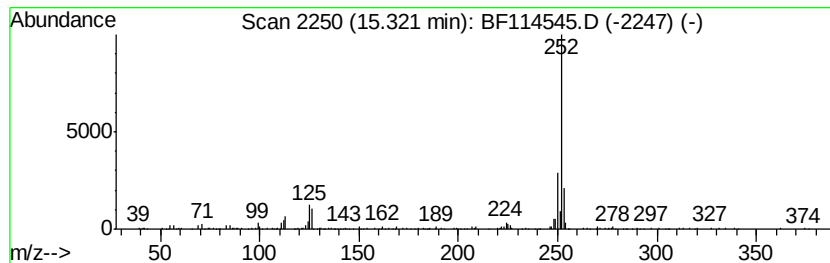
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzo[e]pyrene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.32	2.66 ng	188412	Perylene-d12	15.44

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	98
2			Benzo[a]pyrene	252	C20H12	000050-32-8	96
3			Benzo[k]fluoranthene	252	C20H12	000207-08-9	95
4			Perylene	252	C20H12	000198-55-0	94
5			Benz[e]acephenanthrylene	252	C20H12	000205-99-2	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052319\
Data File : BF114545.D
Acq On : 24 May 2019 1:06
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Sample : K2998-01
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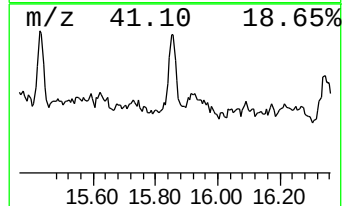
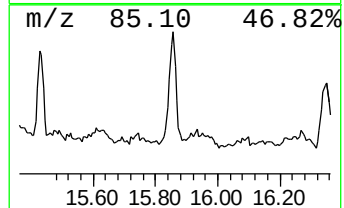
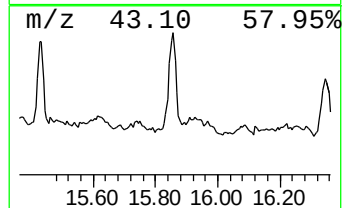
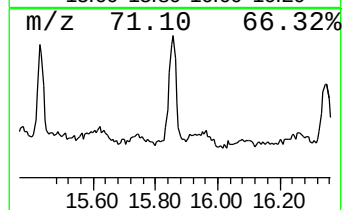
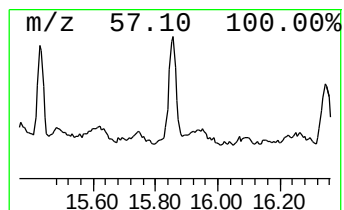
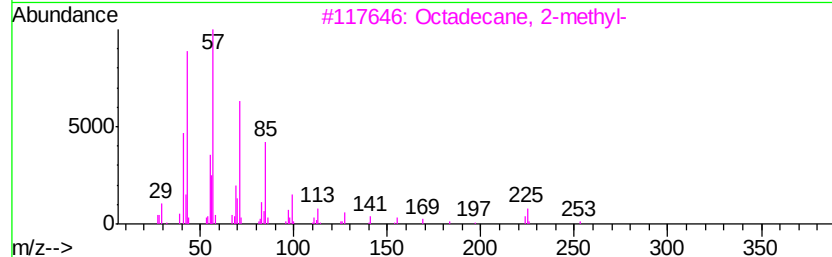
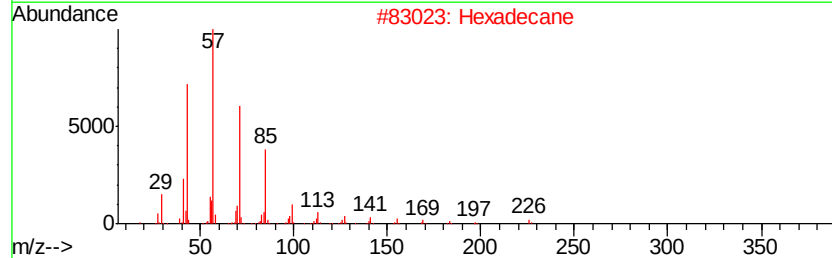
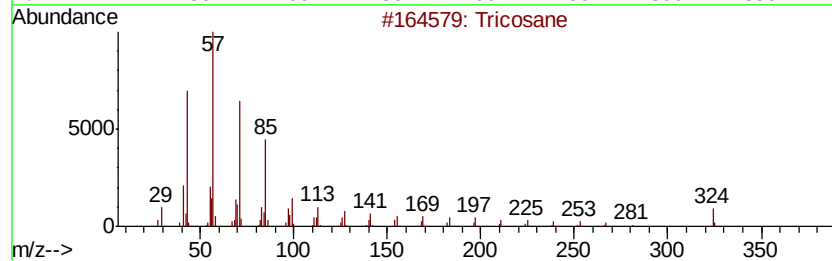
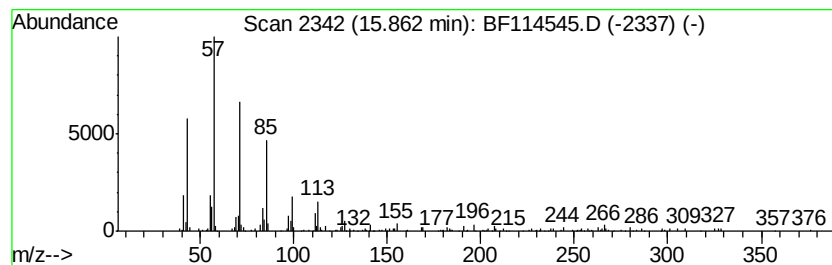
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tricosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.86	4.30 ng	304215	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricosane	324	C23H48	000638-67-5	90
2		Hexadecane	226	C16H34	000544-76-3	86
3		Octadecane, 2-methyl-	268	C19H40	001560-88-9	86
4		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	86
5		2-Bromo dodecane	248	C12H25Br	013187-99-0	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052319\
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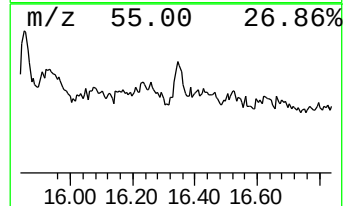
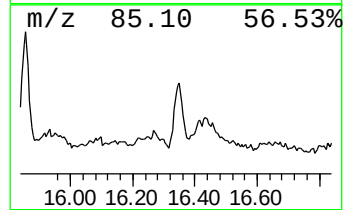
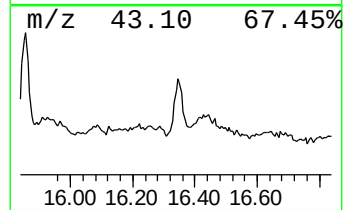
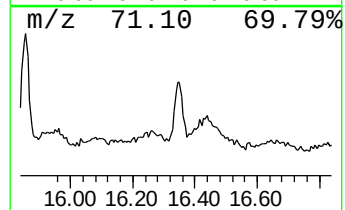
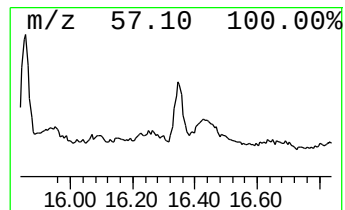
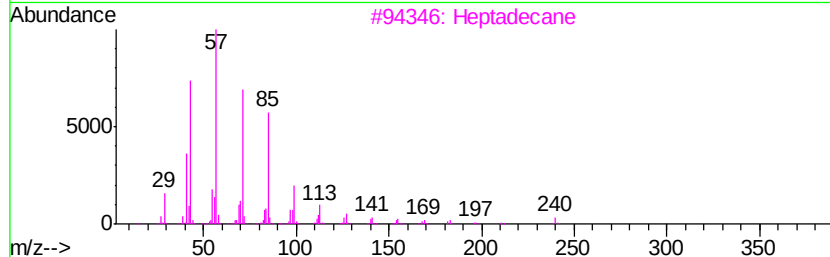
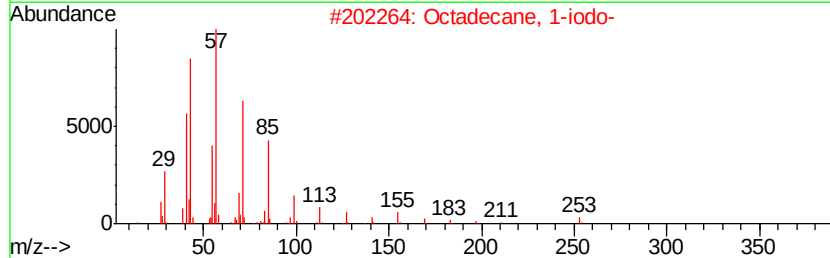
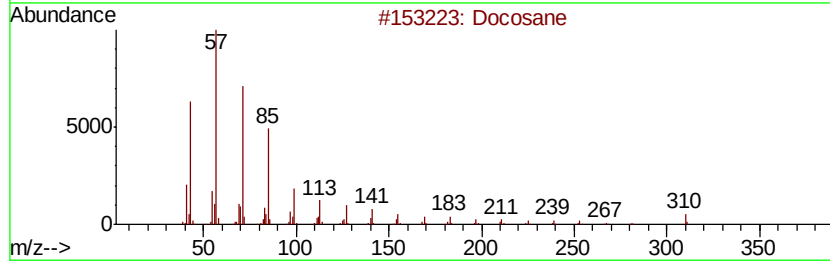
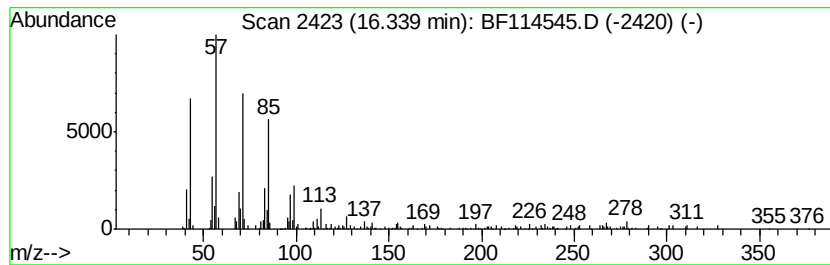
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Docosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.34	3.88 ng	274949	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosane	310	C22H46	000629-97-0	87
2		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	80
3		Heptadecane	240	C17H36	000629-78-7	80
4		Triacontane	422	C30H62	000638-68-6	80
5		Tetracosane	338	C24H50	000646-31-1	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF052319\
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 ALS Vial : 14 Sample Multiplier: 1

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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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n-Hexadecanoic acid	11.86	4.4	ng	359979	4	11.33	1626920	20.0
unknown11.91	11.91	2.1	ng	174851	4	11.33	1626920	20.0
unknown11.93	11.93	6.3	ng	516548	4	11.33	1626920	20.0
18-Norabietane	11.98	51.3	ng	4170490	4	11.33	1626920	20.0
unknown12.08	12.08	51.4	ng	4182550	4	11.33	1626920	20.0
Retene	13.07	120.1	ng	9095570	5	13.98	1514130	20.0
9-Tricosene, (Z)-	13.81	5.6	ng	421466	5	13.98	1514130	20.0
Benzo[e]pyrene	15.32	2.7	ng	188412	6	15.44	1416310	20.0
Tricosane	15.86	4.3	ng	304215	6	15.44	1416310	20.0
Docosane	16.34	3.9	ng	274949	6	15.44	1416310	20.0