

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100319\
 Data File : BF117336.D
 Acq On : 4 Oct 2019 00:11
 Operator : JU
 Sample : K5165-01
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NBP-2019

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-BF091319.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	5.440	567	570	585	rBV	703646	851277	82.66%	7.738%
2	6.463	740	744	761	rBV	653746	968286	94.02%	8.802%
3	6.587	761	765	779	rVB	1094140	1029850	100.00%	9.362%
4	6.810	799	803	810	rBV	290290	265567	25.79%	2.414%
5	6.963	825	829	841	rBV	826740	786582	76.38%	7.150%
6	7.369	895	898	913	rBV	427195	549018	53.31%	4.991%
7	8.092	1017	1021	1043	rBV	203205	279896	27.18%	2.544%
8	9.163	1199	1203	1214	rBV	897311	895335	86.94%	8.139%
9	9.557	1267	1270	1279	rVB	57821	57749	5.61%	0.525%
10	9.704	1290	1295	1300	rBV	9028	10678	1.04%	0.097%
11	9.839	1315	1318	1322	rBV	246675	234677	22.79%	2.133%
12	9.875	1322	1324	1334	rVB	15653	19320	1.88%	0.176%
13	10.633	1449	1453	1464	rBV	371211	398956	38.74%	3.627%
14	10.751	1469	1473	1480	rVB5	7063	13151	1.28%	0.120%
15	11.186	1543	1547	1549	rVB3	15272	13513	1.31%	0.123%
16	11.227	1549	1554	1558	rBV3	7047	10972	1.07%	0.100%
17	11.327	1568	1571	1573	rBV	240528	217395	21.11%	1.976%
18	11.351	1573	1575	1582	rVV	175120	190411	18.49%	1.731%
19	11.404	1582	1584	1591	rVB	31114	37741	3.66%	0.343%
20	11.569	1610	1612	1617	rBV3	12248	14360	1.39%	0.131%
21	11.833	1653	1657	1658	rBV	17568	21489	2.09%	0.195%
22	11.851	1658	1660	1667	rVB3	69880	94052	9.13%	0.855%
23	11.910	1667	1670	1672	rBV3	10503	12171	1.18%	0.111%
24	11.939	1672	1675	1678	rBV2	32780	35251	3.42%	0.320%
25	12.016	1686	1688	1691	rBV4	10376	12455	1.21%	0.113%
26	12.127	1704	1707	1710	rVB	14678	12263	1.19%	0.111%
27	12.163	1710	1713	1717	rBV3	20126	18980	1.84%	0.173%
28	12.274	1726	1732	1735	rBV8	17067	30237	2.94%	0.275%
29	12.380	1747	1750	1752	rBV	26946	22119	2.15%	0.201%
30	12.410	1752	1755	1758	rBV4	15393	16612	1.61%	0.151%
31	12.474	1763	1766	1770	rVB3	21891	27452	2.67%	0.250%
32	12.545	1774	1778	1783	rBV	360399	346601	33.66%	3.151%
33	12.639	1791	1794	1799	rVB6	18893	21106	2.05%	0.192%
34	12.774	1811	1817	1824	rBV	311783	362236	35.17%	3.293%

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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

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35	12.910	1836	1840	1850	rVB	952643	877252	85.18%	7.975%
36	12.992	1850	1854	1858	rBV4	19073	32823	3.19%	0.298%
37	13.168	1882	1884	1886	rBV	17040	14336	1.39%	0.130%
38	13.298	1904	1906	1909	rBV	21356	22811	2.21%	0.207%
39	13.333	1909	1912	1913	rVV	24374	22945	2.23%	0.209%
40	13.374	1917	1919	1922	rVB	21753	20375	1.98%	0.185%
41	13.615	1958	1960	1963	rVB	31453	28422	2.76%	0.258%
42	13.721	1976	1978	1980	rBV	26293	25235	2.45%	0.229%
43	13.745	1980	1982	1985	rVB2	34421	28280	2.75%	0.257%
44	13.774	1985	1987	1988	rBV	21785	18615	1.81%	0.169%
45	13.804	1989	1992	1994	rBV2	96872	92672	9.00%	0.842%
46	13.939	2012	2015	2016	rBV	46160	37114	3.60%	0.337%
47	13.974	2016	2021	2023	rVV2	314759	444361	43.15%	4.039%
48	13.998	2023	2025	2029	rVB	158057	143053	13.89%	1.300%
49	14.421	2095	2097	2099	rBV	49434	51409	4.99%	0.467%
50	15.009	2194	2197	2200	rBV	104165	152368	14.80%	1.385%
51	15.051	2201	2204	2208	rVB	259554	293359	28.49%	2.667%
52	15.309	2245	2248	2253	rVB	60564	70020	6.80%	0.637%
53	15.368	2255	2258	2264	rVB	79172	97180	9.44%	0.883%
54	15.427	2264	2268	2272	rBV2	139590	202106	19.62%	1.837%
55	15.845	2334	2339	2344	rVB	199451	286063	27.78%	2.600%
56	16.898	2513	2518	2527	rVB3	79106	162113	15.74%	1.474%

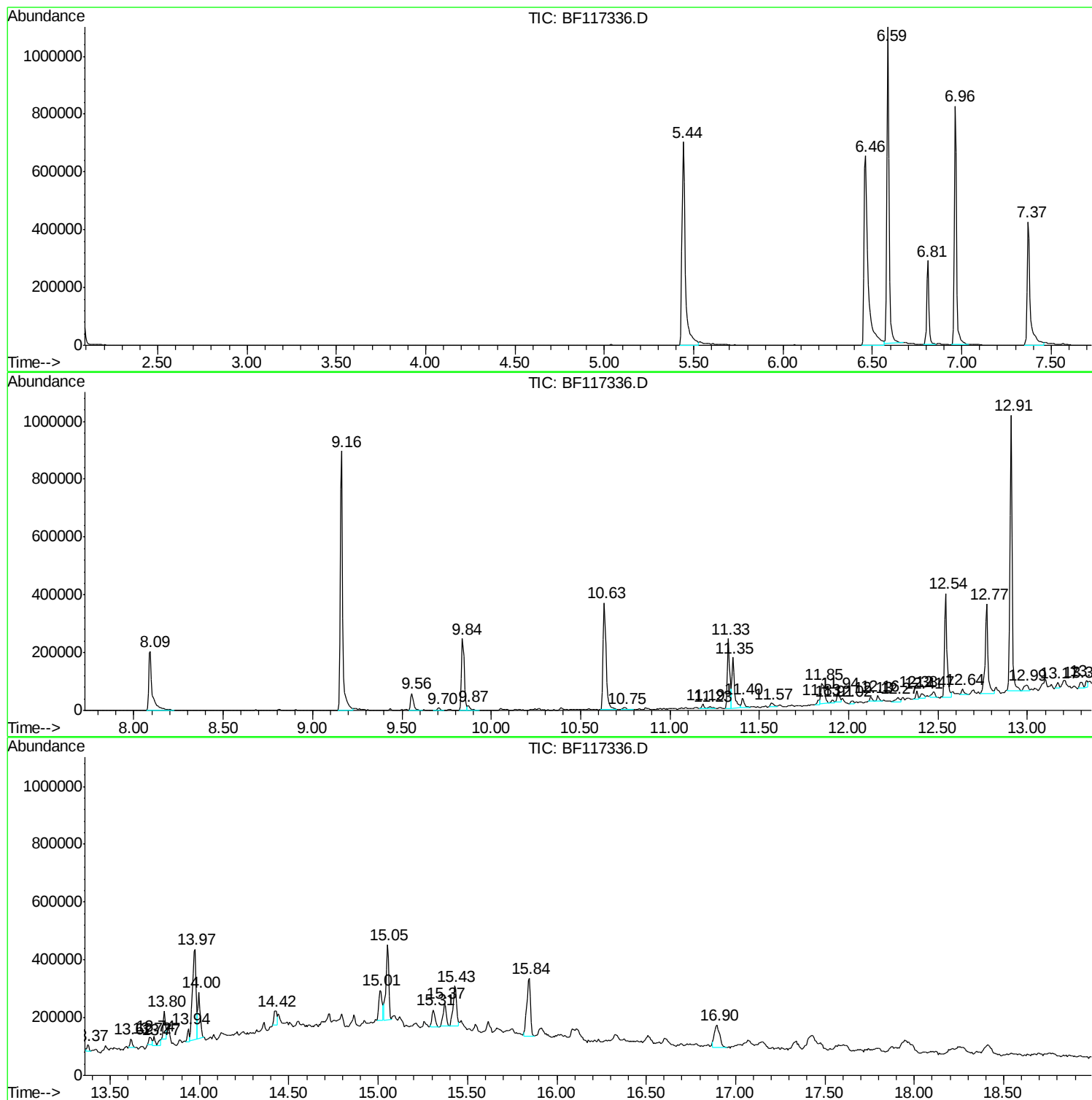
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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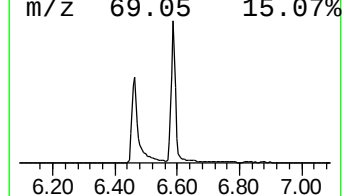
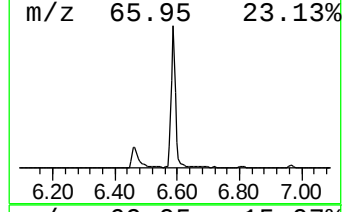
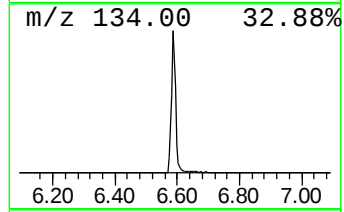
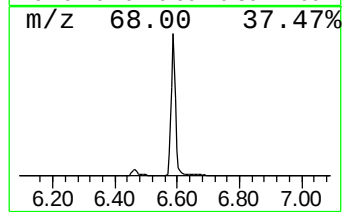
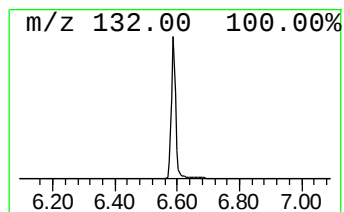
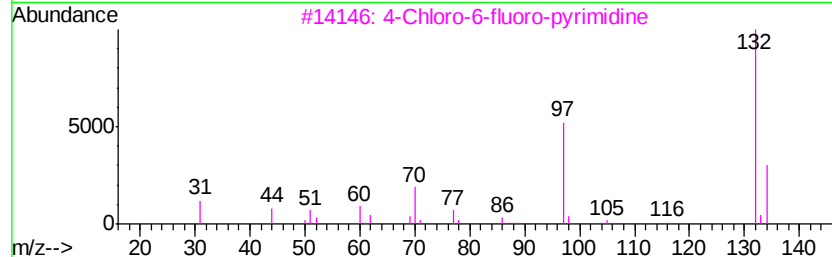
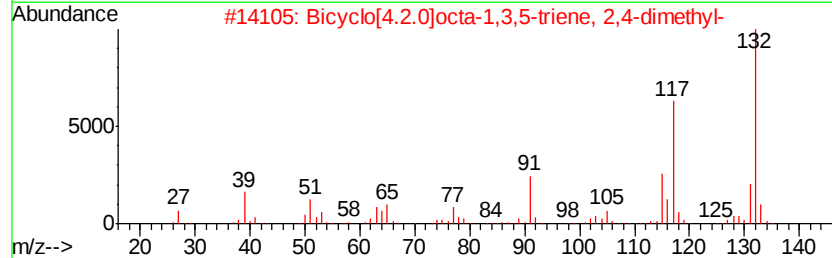
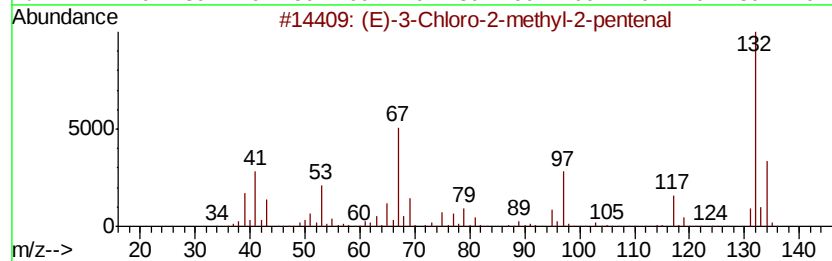
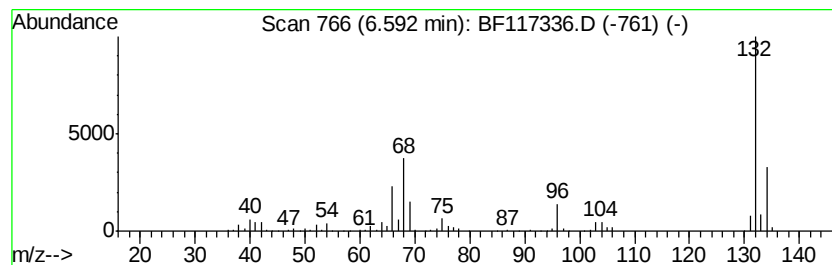
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.59	77.56 ng	1029850	1,4-Dichlorobenzene-d4	6.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
2		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
3		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
4		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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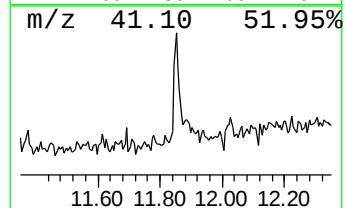
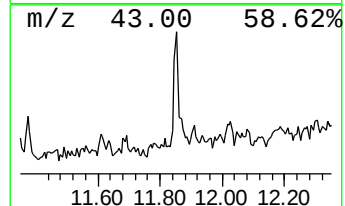
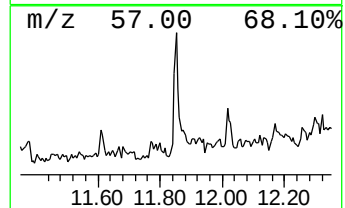
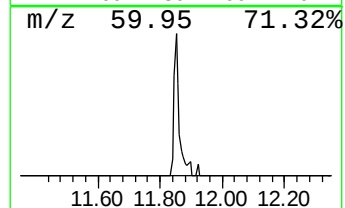
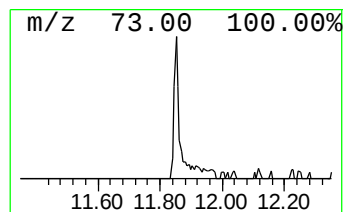
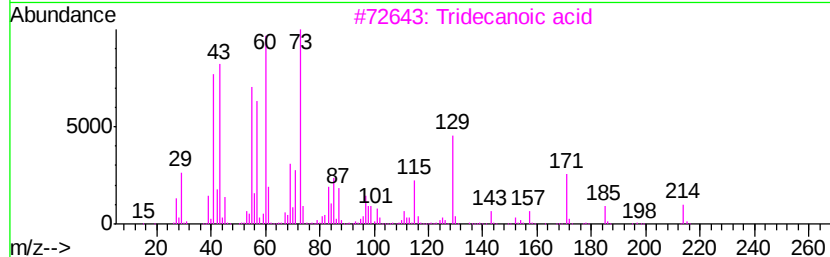
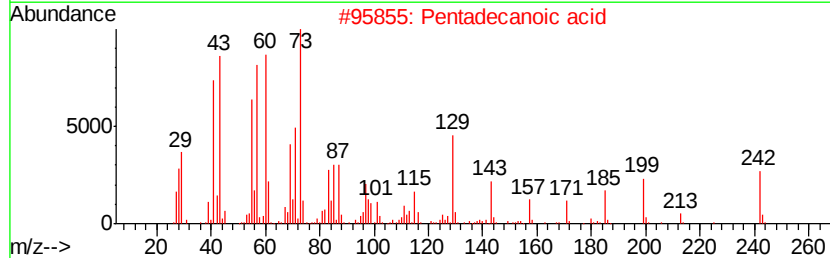
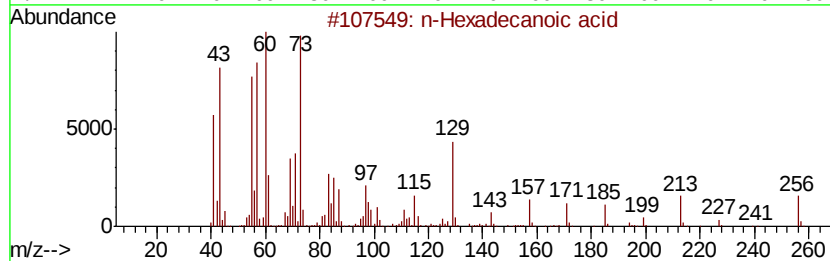
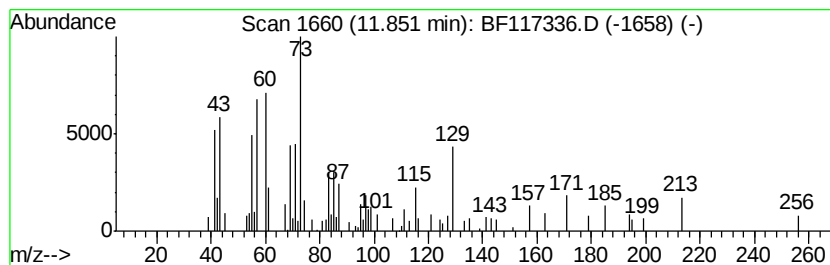
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.85	8.65 ng	94052	Phenanthrene-d10	11.33

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	89
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
3		Tridecanoic acid	214	C13H26O2	000638-53-9	53
4		Lactose	342	C12H22O11	000063-42-3	38
5		2,6,10-Trimethylundeca-1,3-diene	194	C14H26	1000222-09-1	25



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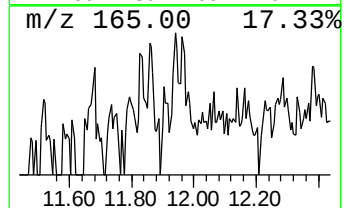
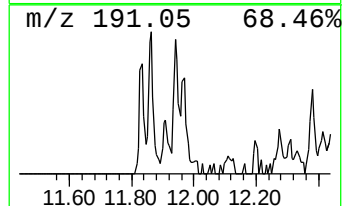
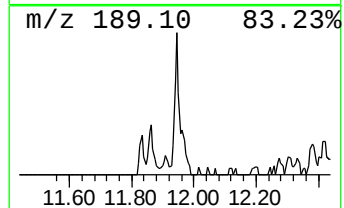
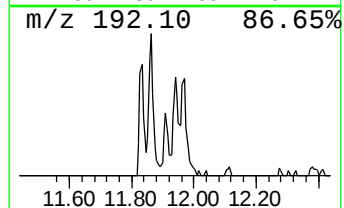
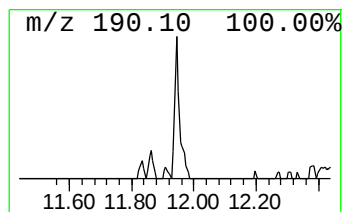
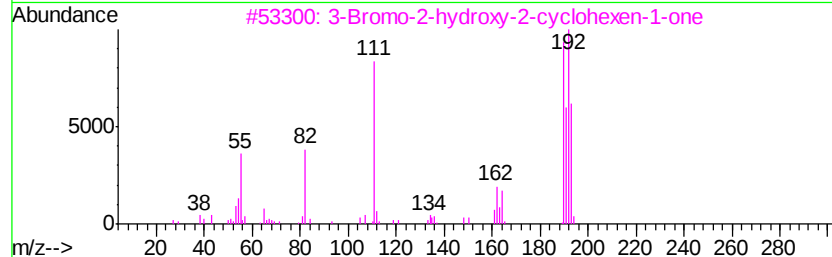
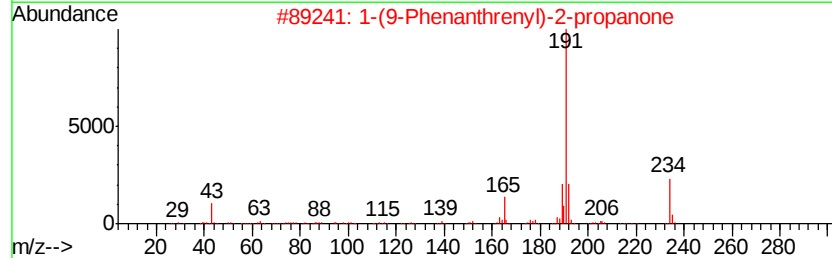
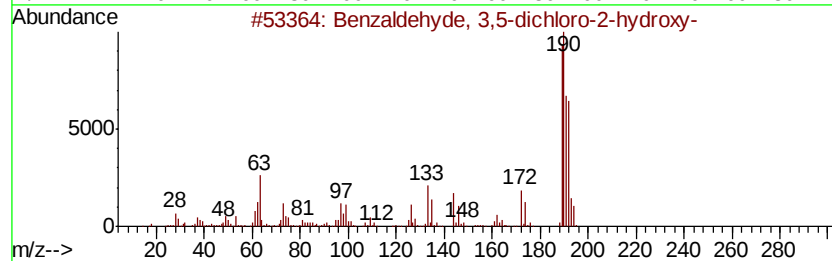
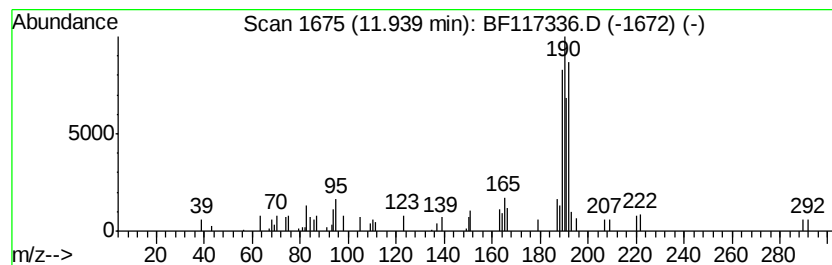
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzaldehyde, 3,5-dichloro-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.94	3.24 ng	35251	Phenanthrene-d10	11.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzaldehydve, 3,5-dichloro-2-hvd...	190	C7H4Cl2O2	000090-60-8	59
2		1-(9-Phenanthrenyl)-2-propanone	234	C17H14O	1000326-08-9	47
3		3-Bromo-2-hydroxy-2-cyclohexen-1...	190	C6H7BrO2	010324-65-9	47
4		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	46
5		3,5-Dichlorobenzoic acid	190	C7H4Cl2O2	000051-36-5	46



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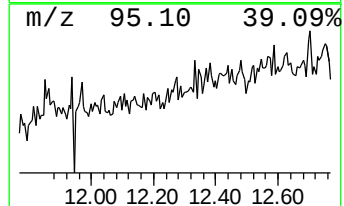
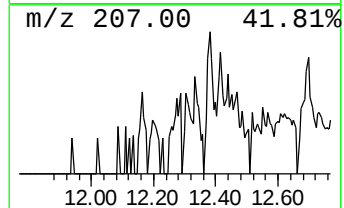
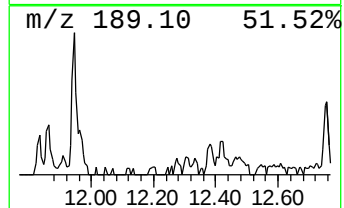
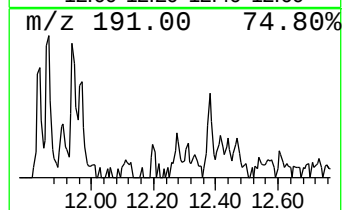
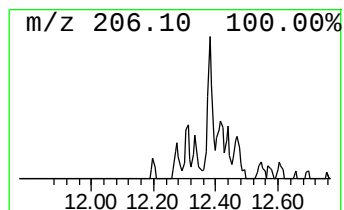
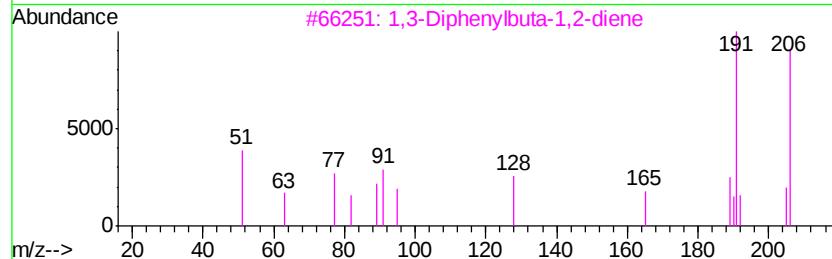
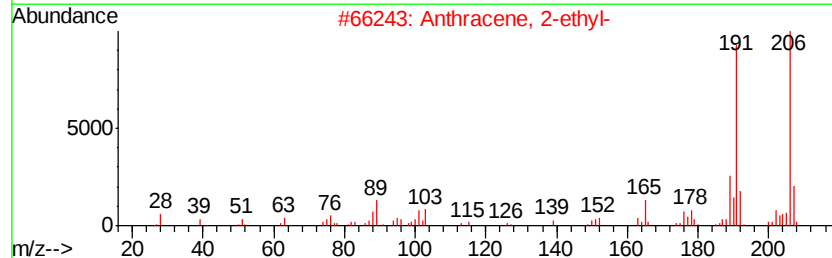
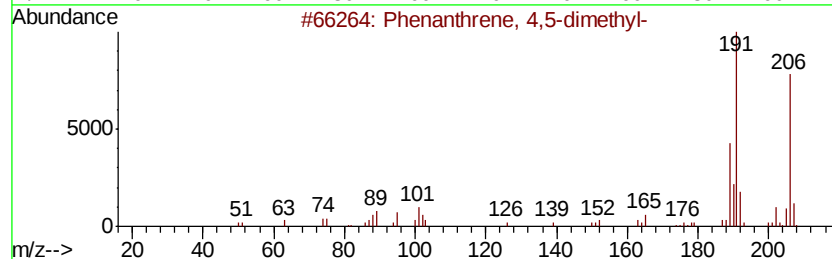
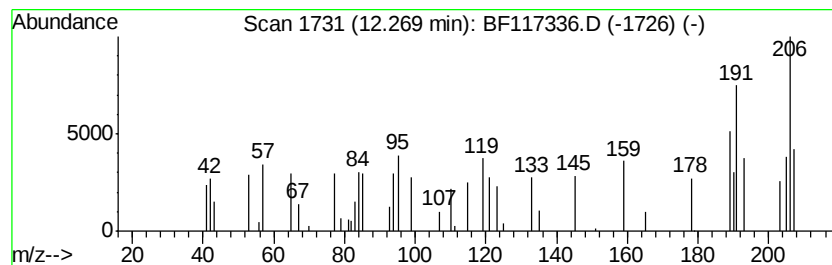
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Peak Number 5 Phenanthrene, 4,5-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.27	2.78 ng	30237	Phenanthrene-d10	11.33

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	53
2		Anthracene, 2-ethyl-	206	C16H14	052251-71-5	53
3		1,3-Diphenylbuta-1,2-diene	206	C16H14	053544-89-1	46
4		3,4-Dimethoxybenzylideneacetone	206	C12H14O3	015001-27-1	46
5		1,4-Diphenyl-1,3-butadiene	206	C16H14	000886-65-7	43



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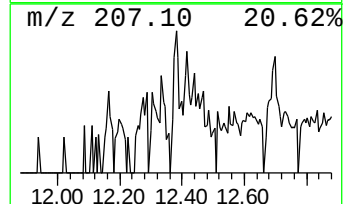
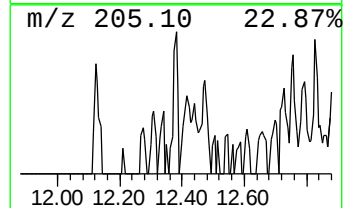
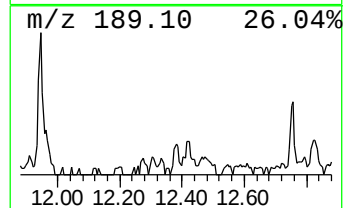
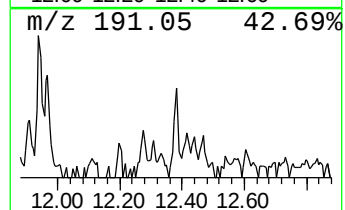
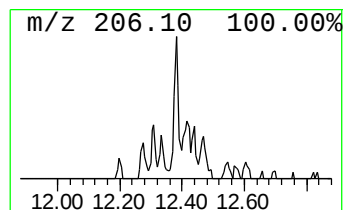
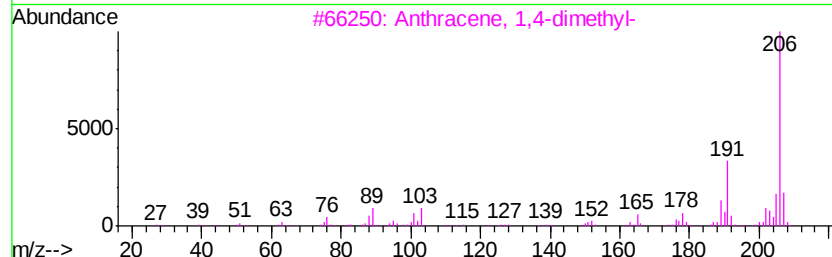
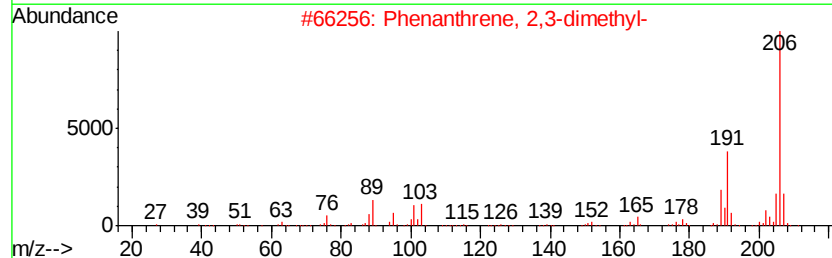
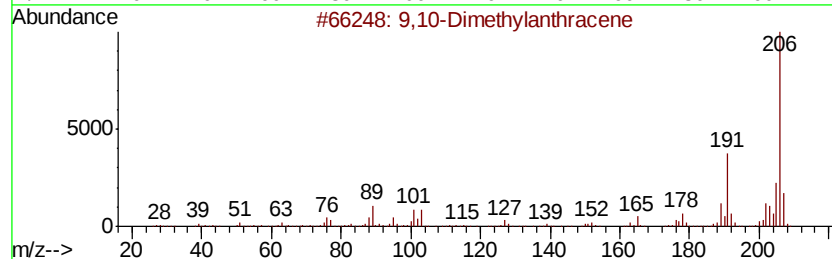
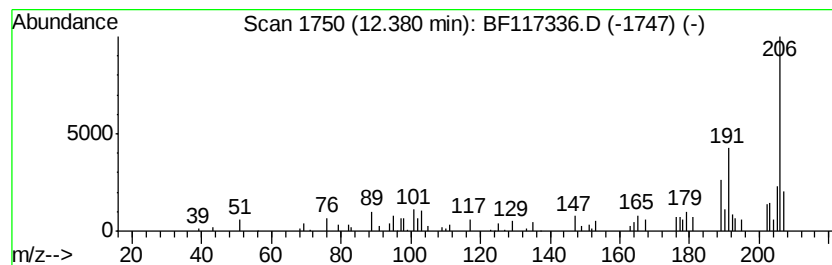
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 9,10-Dimethylantracene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.38	2.03 ng	22119	Phenanthrene-d10	11.33

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,10-Dimethylantracene	206	C16H14	000781-43-1	96
2			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	93
3			Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	91
4			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	90
5			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	89



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100319\
Data File : BF117336.D
Acq On : 4 Oct 2019 00:11
Operator : JU
Sample : K5165-01
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
NBP-2019

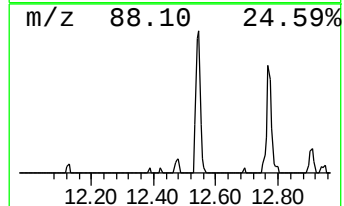
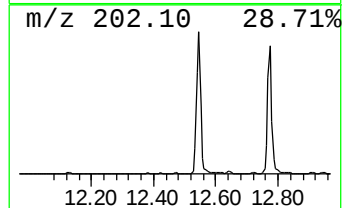
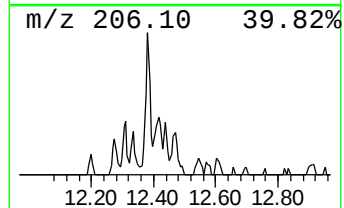
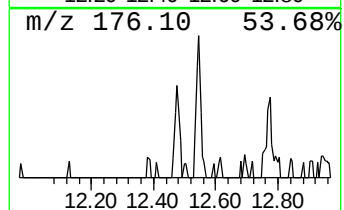
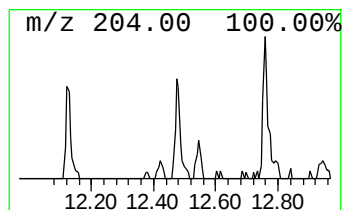
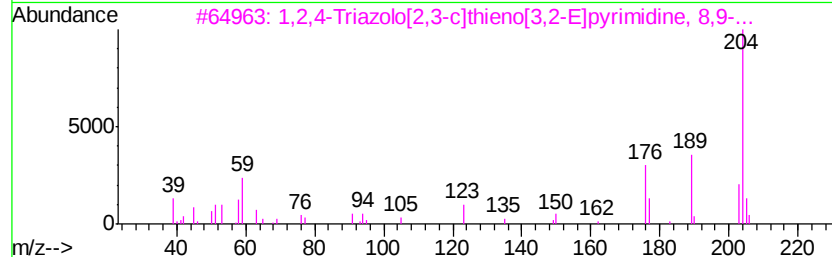
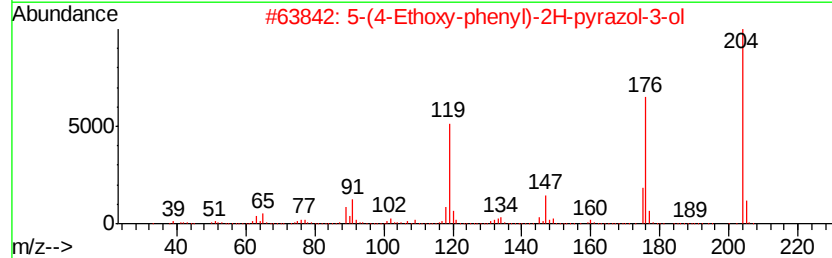
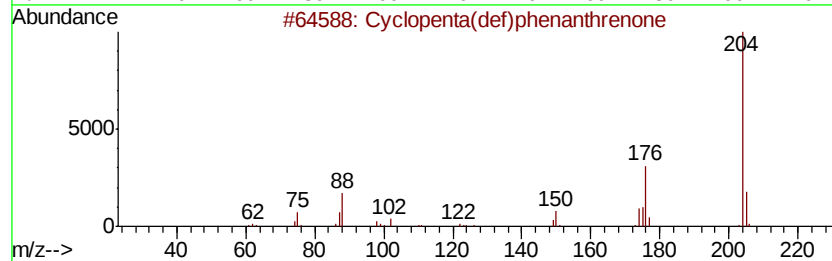
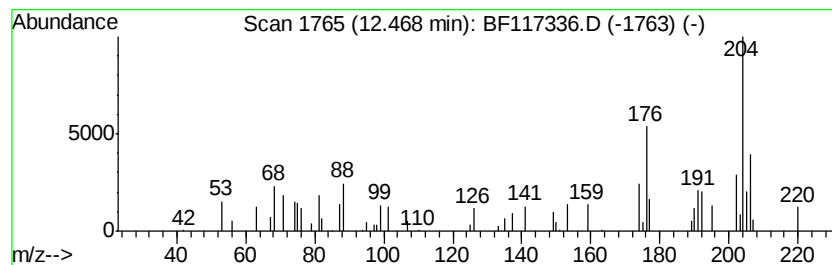
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Cyclopenta(def)phenanthrenone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.47	2.53 ng	27452	Phenanthrene-d10	11.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	68
2		5-(4-Ethoxy-phenyl)-2H-pyrazol-3-ol	204	C11H12N2O2	1000278-26-8	53
3		1,2,4-Triazolo[2,3-c]thieno[3,2-...	204	C9H8N4S	113246-88-1	53
4		[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	47
5		4-Methoxy-2-(4-fluorophenyl)pyri...	204	C11H9FN2O	076128-75-1	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100319\
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Operator : JU
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Misc :
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Instrument :
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ClientSampleId :
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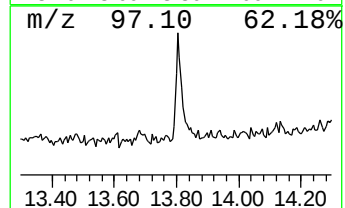
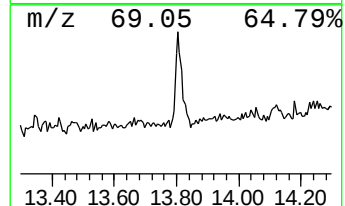
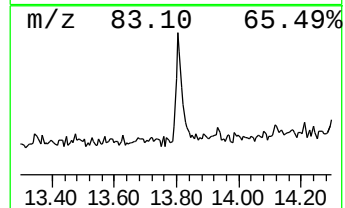
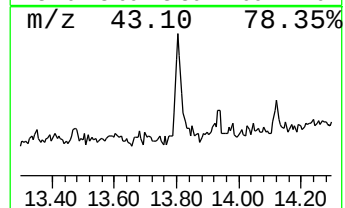
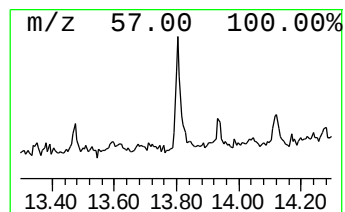
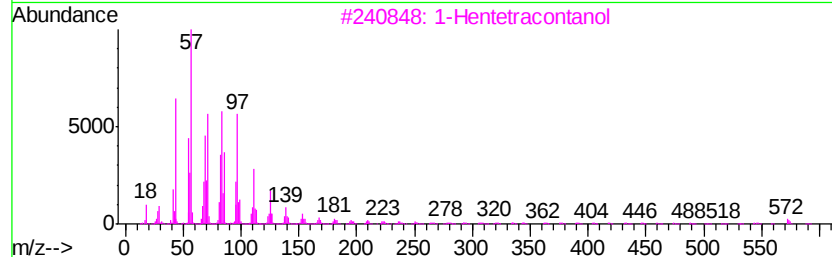
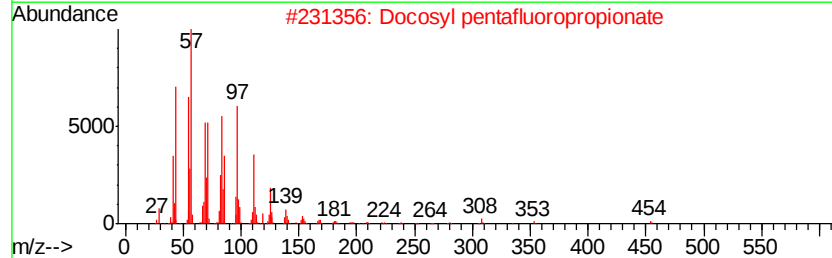
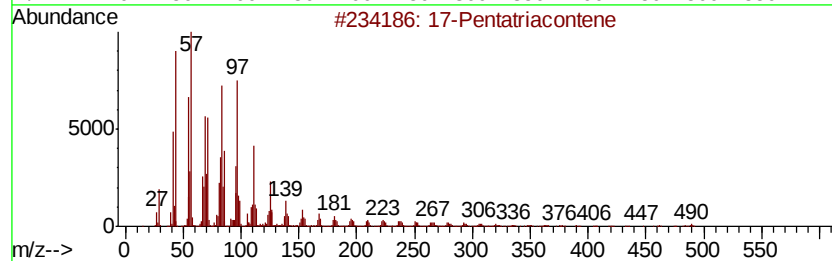
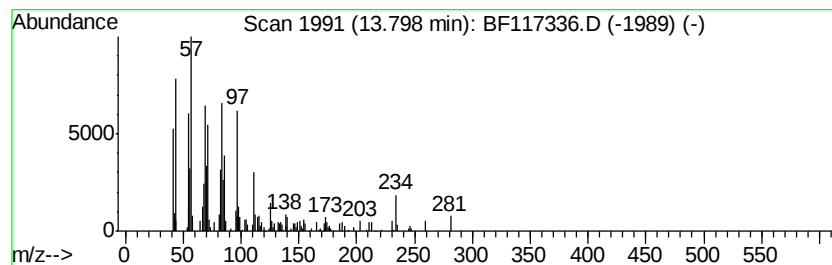
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 17-Pentatriacontene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.80	4.17 ng	92672	Chrysene-d12	13.97

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		17-Pentatriacontene	491	C35H70	006971-40-0	87
2		Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	86
3		1-Hentetracontanol	593	C41H84O	040710-42-7	83
4		Isobutyl tetradecyl carbonate	314	C19H38O3	959275-58-2	80
5		Tetracosyl heptafluorobutyrate	550	C28H49F7O2	1000351-83-7	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100319\
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Misc :
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ClientSampleId :
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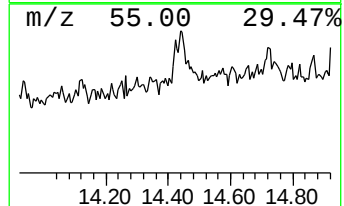
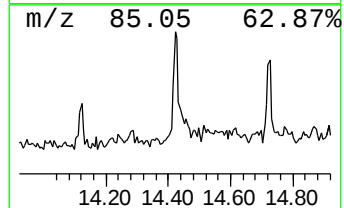
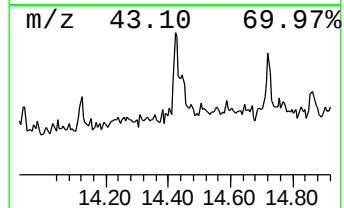
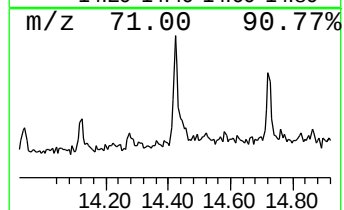
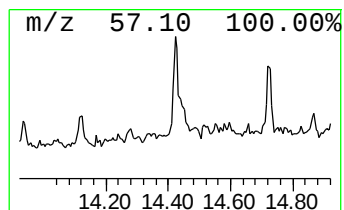
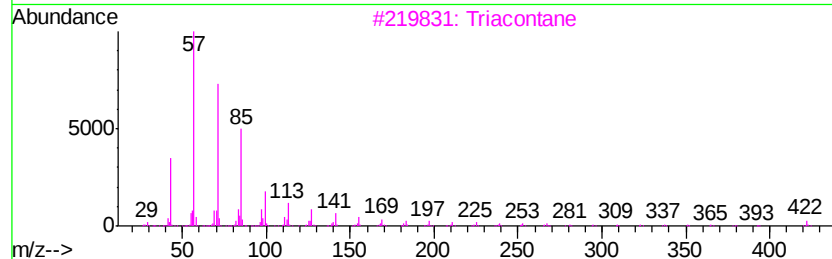
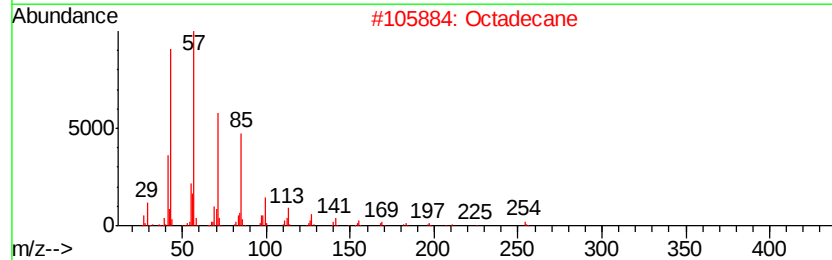
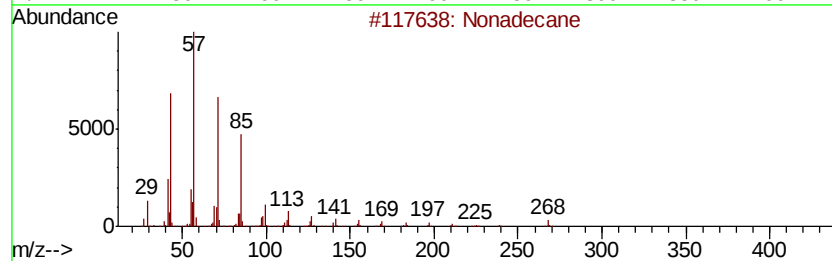
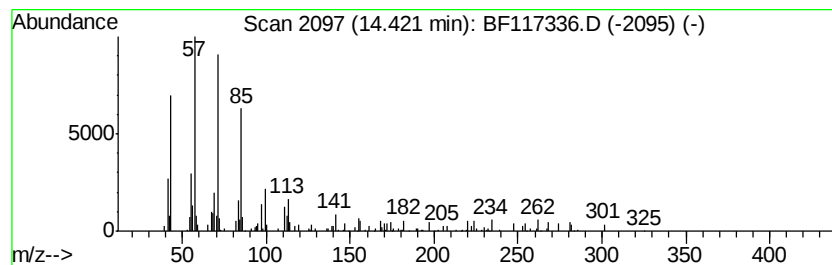
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Nonadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.42	2.31 ng	51409	Chrysene-d12	13.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	81
2		Octadecane	254	C18H38	000593-45-3	74
3		Triacotane	422	C30H62	000638-68-6	72
4		Octacosane	394	C28H58	000630-02-4	72
5		Eicosane	282	C20H42	000112-95-8	68



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100319\
Data File : BF117336.D
Acq On : 4 Oct 2019 00:11
Operator : JU
Sample : K5165-01
Misc :
ALS Vial : 22 Sample Multiplier: 1

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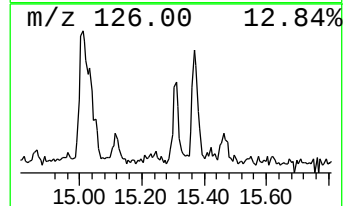
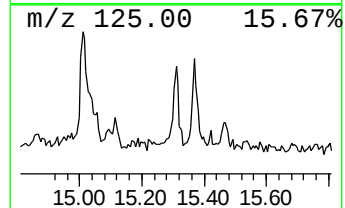
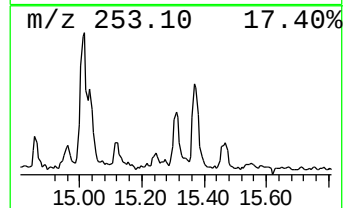
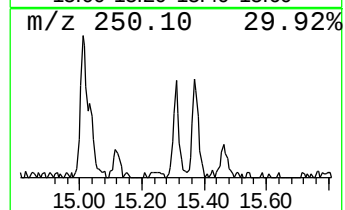
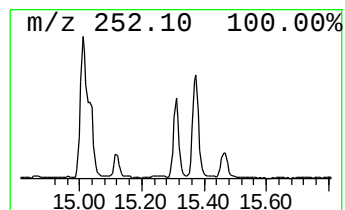
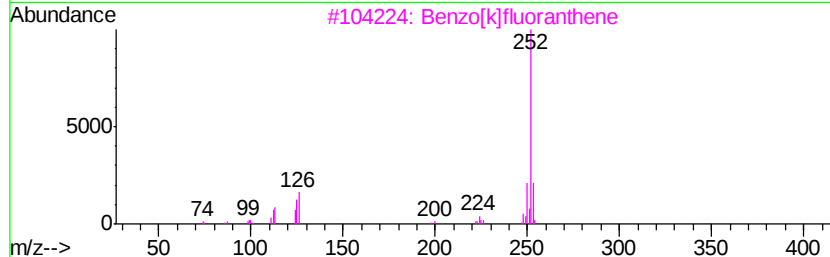
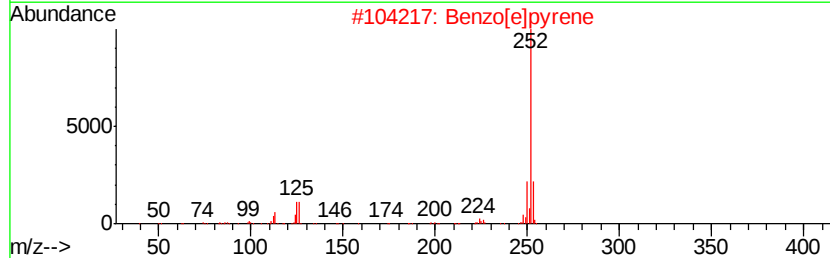
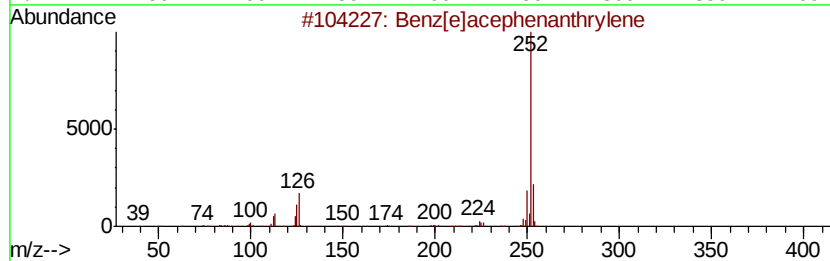
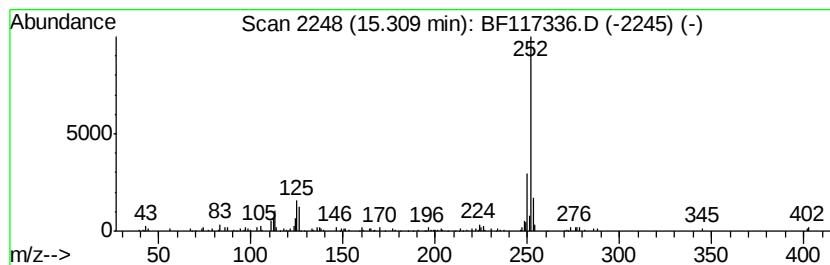
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzo[e]pyrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.31	6.93 ng	70020	Perylene-d12	15.43

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	97
2	Benzo[e]pyrene	252	C20H12	000192-97-2	96
3	Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
4	Benzo[a]pyrene	252	C20H12	000050-32-8	89
5	Perylene	252	C20H12	000198-55-0	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100319\
 Data File : BF117336.D
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Instrument :
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 ClientSampleId :
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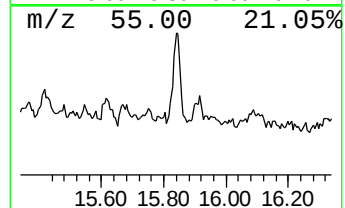
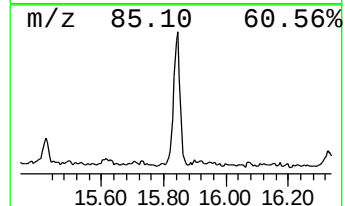
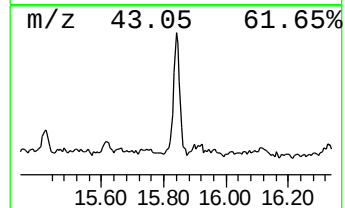
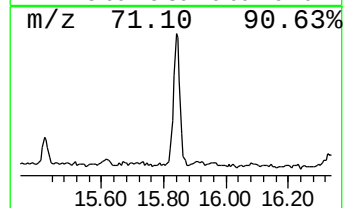
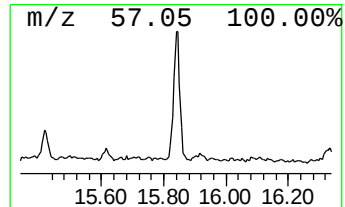
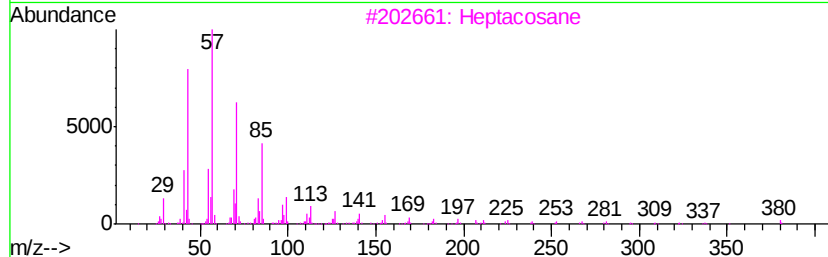
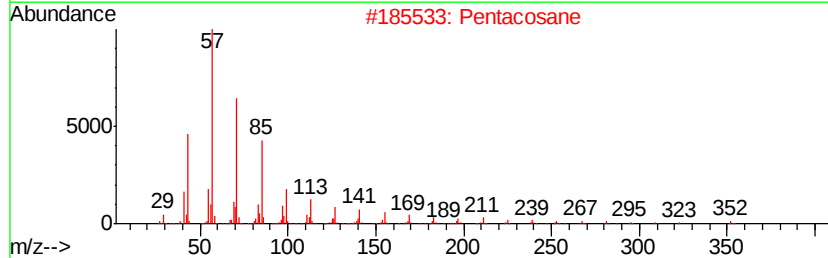
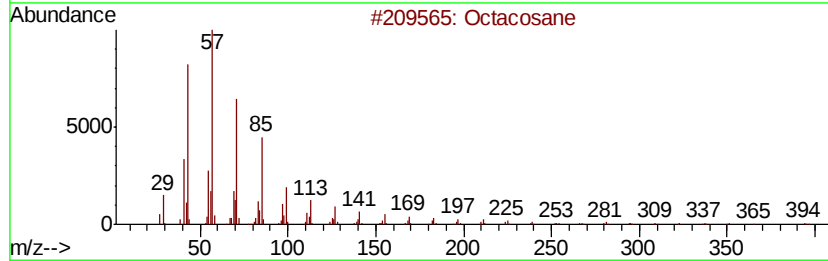
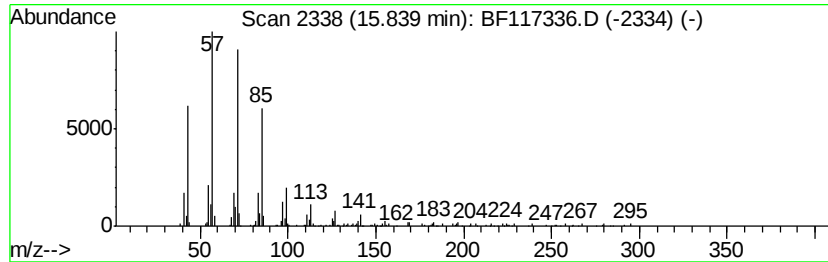
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 11 Octacosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.84	28.31 ng	286063	Perylene-d12	15.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Pentacosane	352	C25H52	000629-99-2	90
3		Heptacosane	380	C27H56	000593-49-7	90
4		2-methyloctacosane	408	C29H60	1000376-72-8	90
5		Hentriacontane	437	C31H64	000630-04-6	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF100319\
Data File : BF117336.D
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Misc :
ALS Vial : 22 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	11.85	8.7	ng	94052	4	11.33	217395	20.0
Benzaldehyde, 3,5...	11.94	3.2	ng	35251	4	11.33	217395	20.0
Phenanthrene, 4,5...	12.27	2.8	ng	30237	4	11.33	217395	20.0
9,10-Dimethylanthe...	12.38	2.0	ng	22119	4	11.33	217395	20.0
Cyclopenta(def)ph...	12.47	2.5	ng	27452	4	11.33	217395	20.0
17-Pentatriacontene	13.80	4.2	ng	92672	5	13.97	444361	20.0
Nonadecane	14.42	2.3	ng	51409	5	13.97	444361	20.0
Benzo[e]pyrene	15.31	6.9	ng	70020	6	15.43	202106	20.0
Octacosane	15.84	28.3	ng	286063	6	15.43	202106	20.0