

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042419\
 Data File : BF113951.D
 Acq On : 24 Apr 2019 10:46
 Operator : JU/SJ
 Sample : K2522-07
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-WOODBURY

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-BF042219.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	2.210	17	21	47	rVB2	74218	231984	5.81%	0.659%
2	5.522	579	584	587	rBV	2909395	2966978	74.31%	8.427%
3	6.528	749	755	766	rBV	2661776	3105963	77.79%	8.821%
4	6.669	774	779	782	rBV	3360598	3191405	79.93%	9.064%
5	6.892	814	817	821	rBV	819245	739910	18.53%	2.101%
6	7.051	840	844	848	rBV	2814451	2558805	64.08%	7.267%
7	7.451	907	912	915	rBV	1937020	1973663	49.43%	5.605%
8	8.175	1031	1035	1050	rVB	1092682	947169	23.72%	2.690%
9	9.251	1213	1218	1221	rBV	4115645	3783397	94.75%	10.745%
10	9.375	1235	1239	1250	rVB	127191	166070	4.16%	0.472%
11	9.639	1281	1284	1289	rBV	381014	309496	7.75%	0.879%
12	9.927	1329	1333	1337	rBV	1151765	1129023	28.28%	3.207%
13	10.716	1463	1467	1481	rBV	2288620	2167930	54.29%	6.157%
14	11.416	1582	1586	1588	rBV	1432626	1225313	30.69%	3.480%
15	11.433	1588	1589	1592	rVB	82333	63706	1.60%	0.181%
16	11.945	1668	1676	1687	rBV	700490	966832	24.21%	2.746%
17	12.627	1783	1792	1795	rBV2	261055	582159	14.58%	1.653%
18	12.657	1795	1797	1804	rVV2	164804	207090	5.19%	0.588%
19	12.727	1804	1809	1821	rVV	642355	786762	19.70%	2.234%
20	12.851	1826	1830	1833	rBV	106641	117638	2.95%	0.334%
21	12.998	1851	1855	1859	rBV	4080957	3992906	100.00%	11.340%
22	13.227	1891	1894	1900	rBV2	86717	100536	2.52%	0.286%
23	13.533	1944	1946	1949	rVB	81547	65676	1.64%	0.187%
24	13.574	1949	1953	1956	rVB	126711	116448	2.92%	0.331%
25	13.898	2005	2008	2015	rVB	170386	166414	4.17%	0.473%
26	14.051	2029	2034	2037	rBV	1210380	1175550	29.44%	3.339%
27	14.215	2059	2062	2068	rBV	139986	128363	3.21%	0.365%
28	14.527	2111	2115	2119	rBV	136419	133879	3.35%	0.380%
29	14.839	2164	2168	2177	rVB2	137506	142816	3.58%	0.406%
30	15.104	2209	2213	2216	rBV2	44169	66733	1.67%	0.190%
31	15.186	2222	2227	2230	rBV	137388	164701	4.12%	0.468%
32	15.468	2272	2275	2280	rVB	35335	46383	1.16%	0.132%
33	15.539	2281	2287	2290	rVV	682230	903525	22.63%	2.566%
34	15.568	2290	2292	2299	rVB2	154650	197435	4.94%	0.561%

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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	16.015	2363	2368	2374	rVB	114226	168786	4.23%	0.479%
36	16.533	2451	2456	2462	rVB	64490	110336	2.76%	0.313%
37	17.139	2554	2559	2564	rBV2	26049	49825	1.25%	0.142%
38	18.615	2801	2810	2821	rBV4	78141	258378	6.47%	0.734%

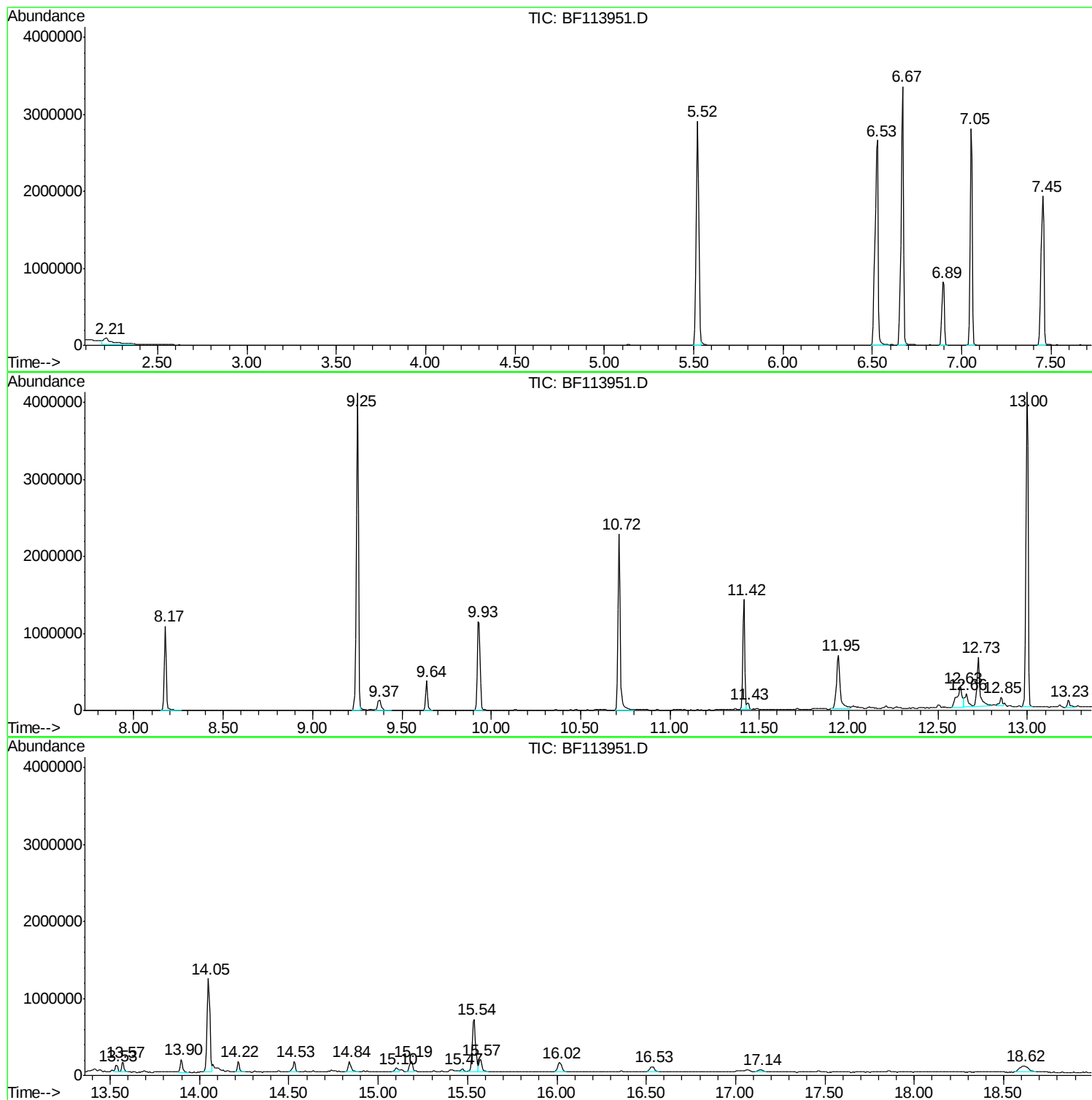
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF042219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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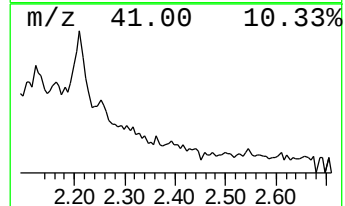
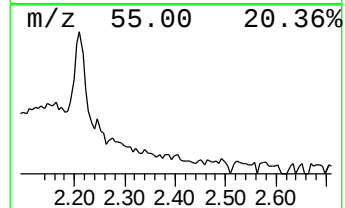
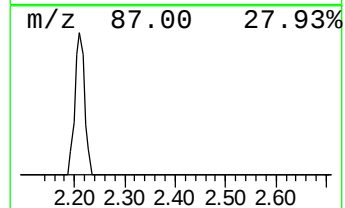
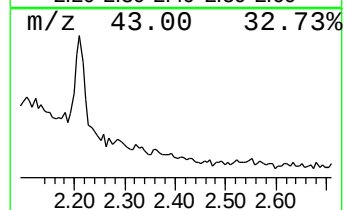
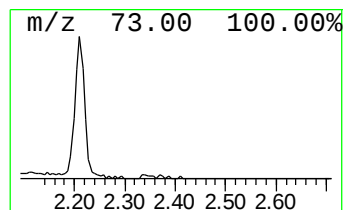
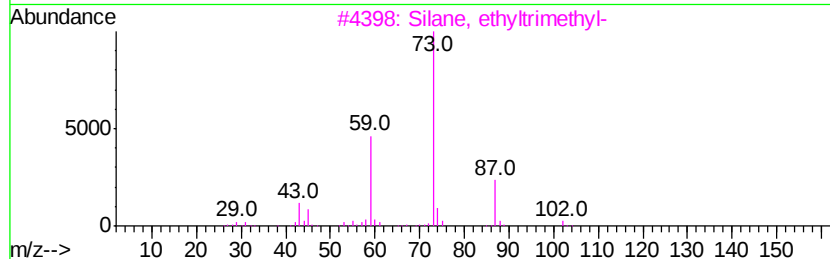
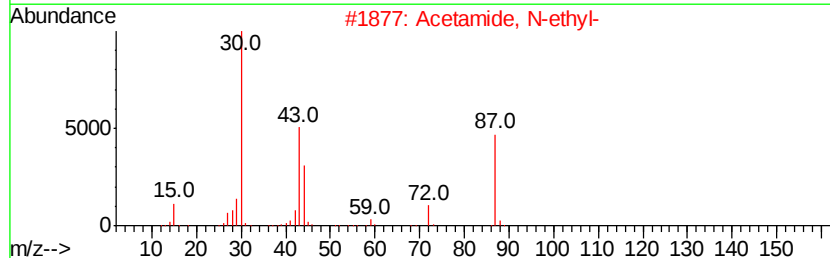
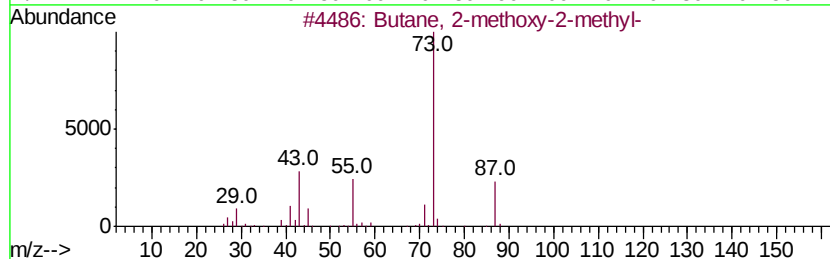
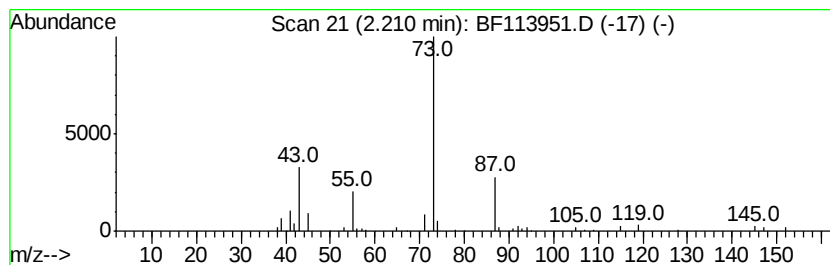
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF042219.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.21	6.27 ng	231984	1,4-Dichlorobenzene-d4	6.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	59
2		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
3		Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	9
4		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	9
5		1-(2-Methoxyethoxy)-2-methyl-2-p...	162	C8H18O3	1000364-24-4	9



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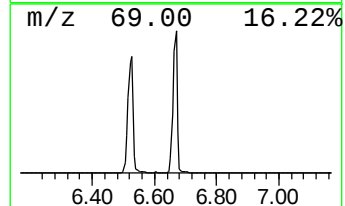
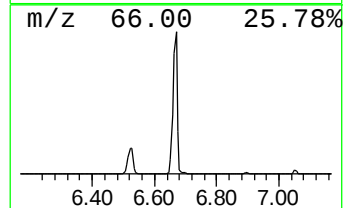
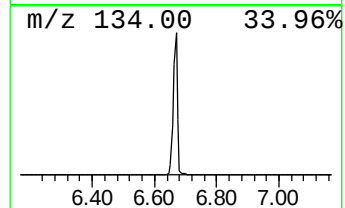
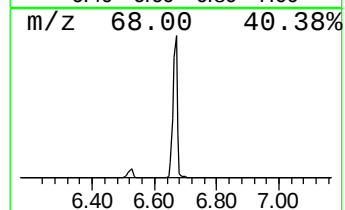
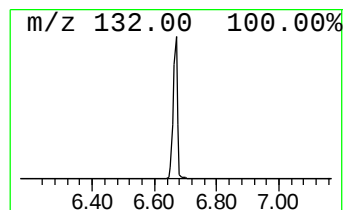
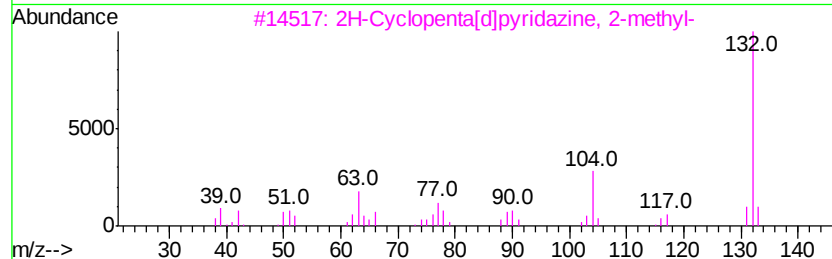
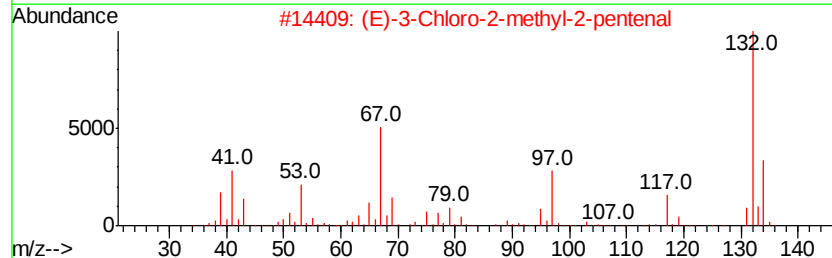
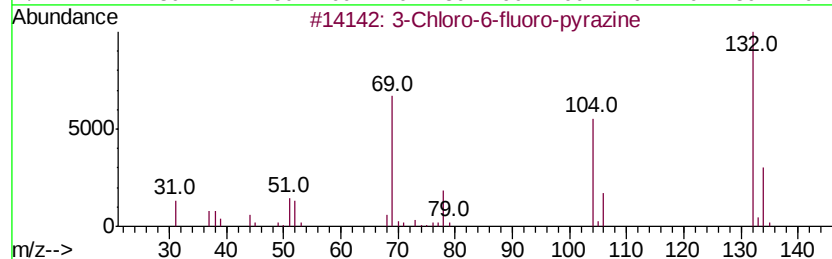
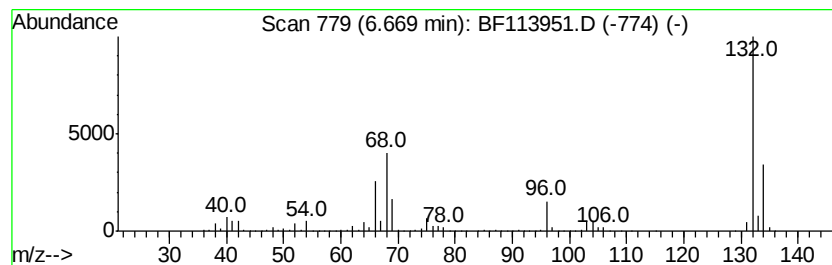
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 unknown6.67 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.67	86.26 ng	3191410	1,4-Dichlorobenzene-d4	6.90

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		2H-Cyclopentaf[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
4		Xenon	132	Xe	007440-63-3	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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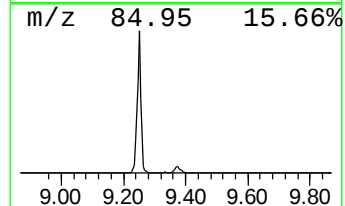
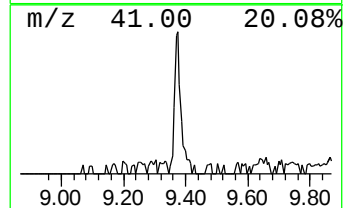
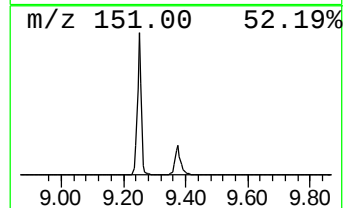
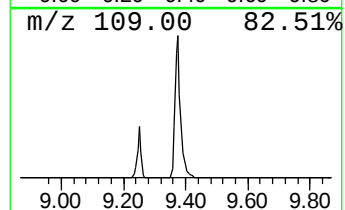
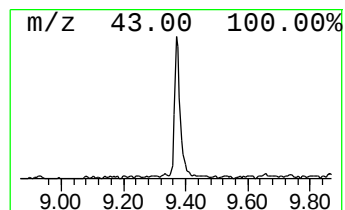
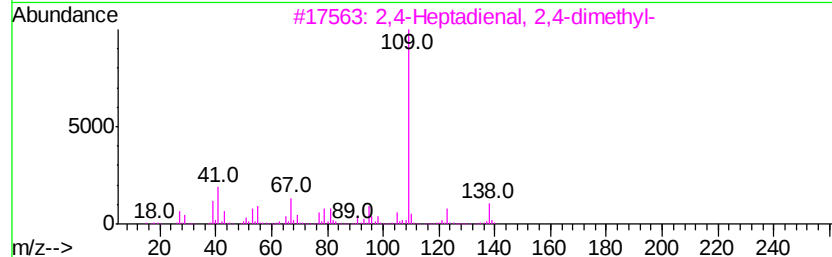
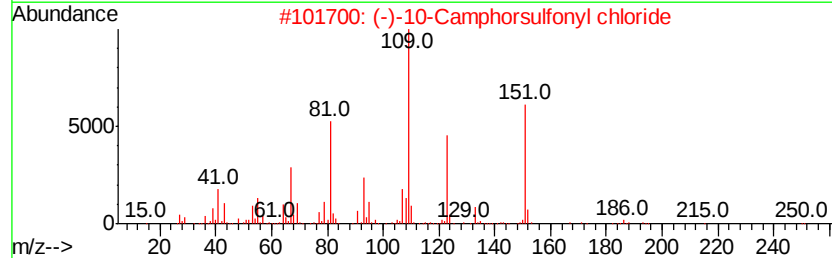
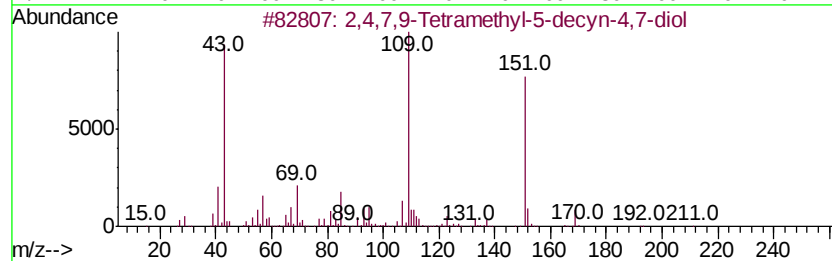
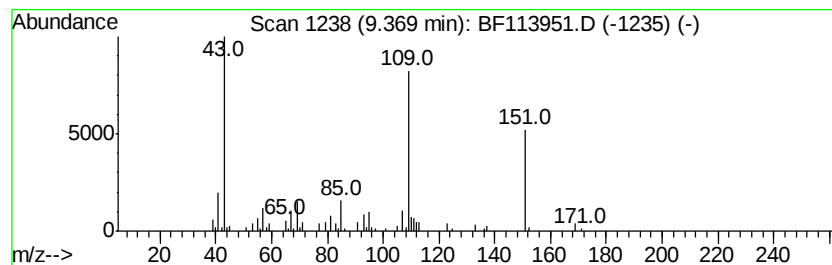
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.37	2.94 ng	166070	Acenaphthene-d10	9.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	91
2		(-)-10-Camphorsulfonyl chloride	250	C10H15ClO3S	039262-22-1	50
3		2,4-Heptadienal, 2,4-dimethyl-	138	C9H14O	042452-48-2	43
4		Metacetamol	151	C8H9NO2	000621-42-1	40
5		1-(1,2,3-Trimethyl-cyclopent-2-e...	152	C10H16O	070987-81-4	38



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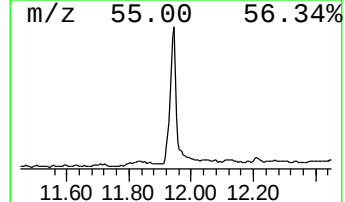
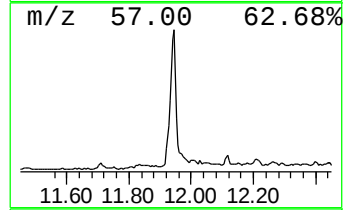
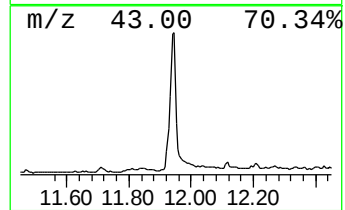
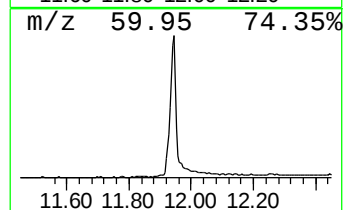
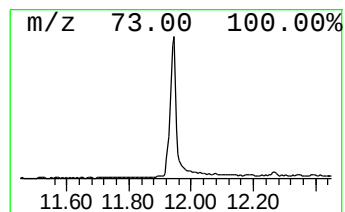
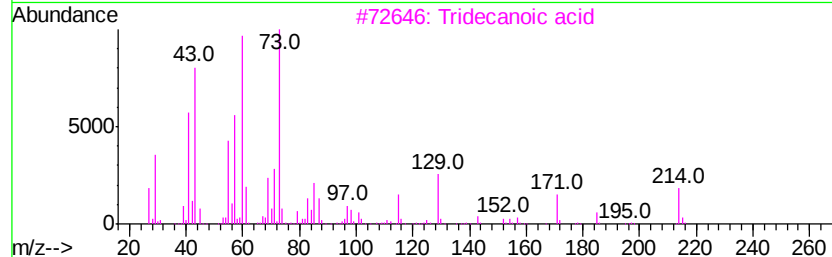
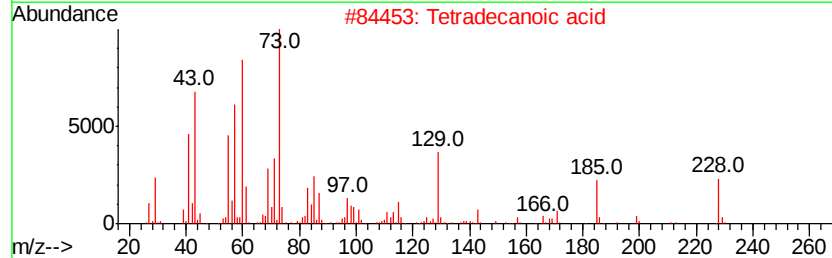
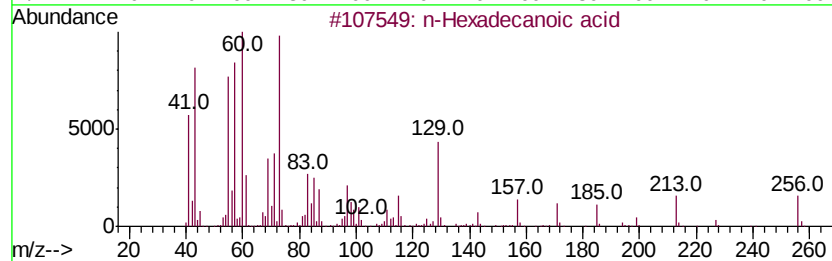
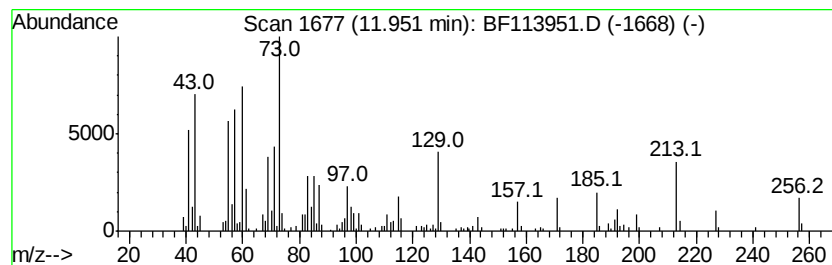
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Peak Number 5 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.95	15.78 ng	966832	Phenanthrene-d10	11.42

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	94
3		Tridecanoic acid	214	C13H26O2	000638-53-9	91
4		n-Decanoic acid	172	C10H20O2	000334-48-5	70
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	68



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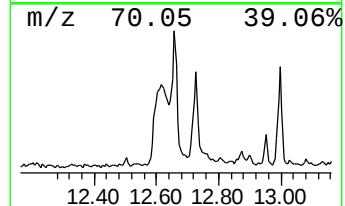
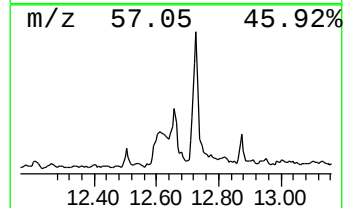
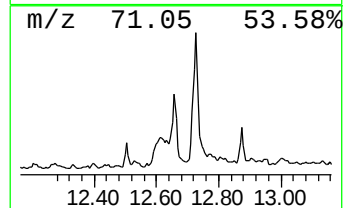
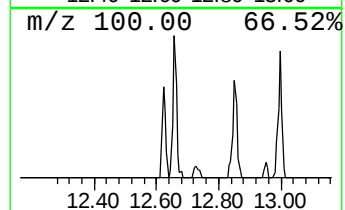
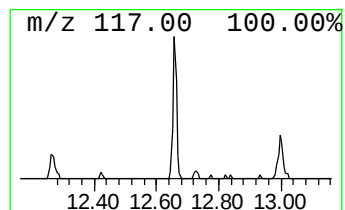
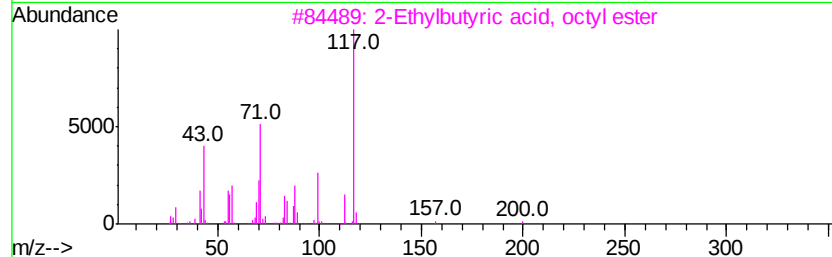
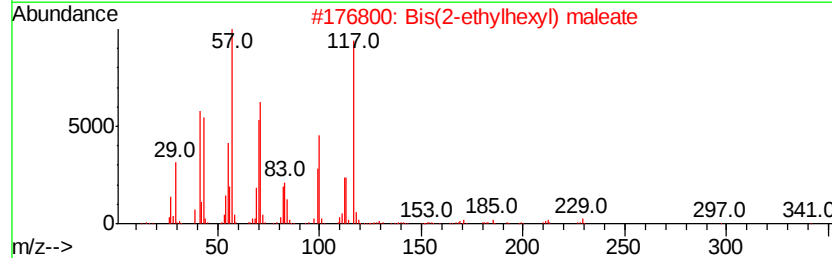
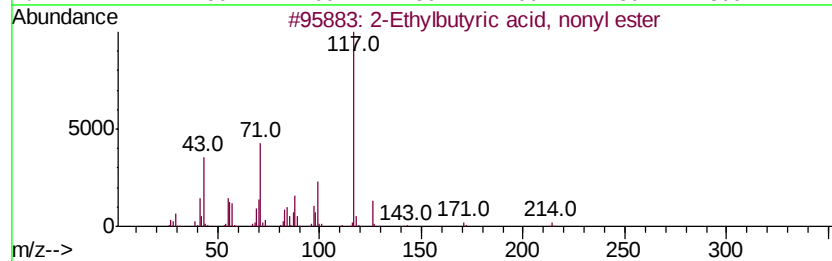
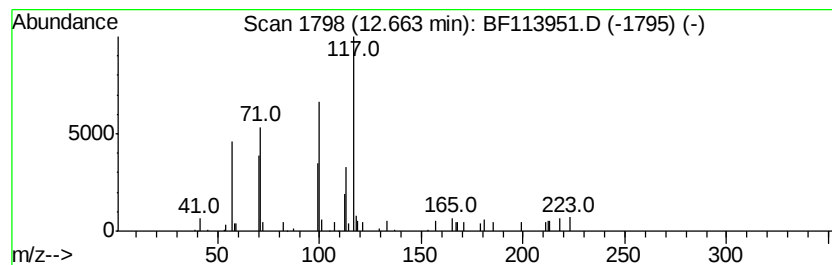
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 2-Ethylbutyric acid, nonyl ... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.66	3.38 ng	207090	Phenanthrene-d10	11.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Ethylbutyric acid, nonyl ester	242	C15H30O2	1000369-43-9	53
2		Bis(2-ethylhexyl) maleate	340	C20H36O4	000142-16-5	50
3		2-Ethylbutyric acid, octyl ester	228	C14H28O2	1000340-24-3	42
4		Carbonic acid, neopentyl 2-ethyl...	244	C14H28O3	1000357-85-7	38
5		Oxalic acid, 2-ethylhexyl octyl ...	314	C18H34O4	1000309-39-1	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042419\
Data File : BF113951.D
Acq On : 24 Apr 2019 10:46
Operator : JU/SJ
Sample : K2522-07
Misc :
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Instrument :
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ClientSampleId :
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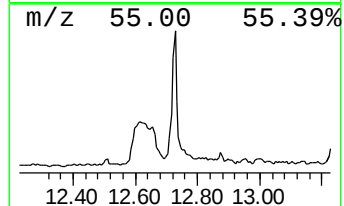
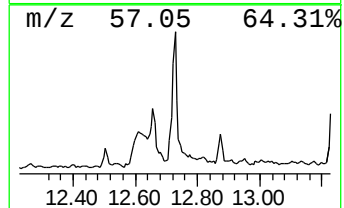
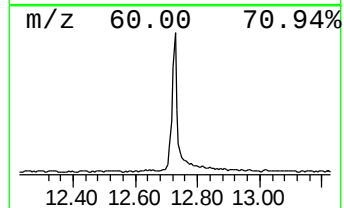
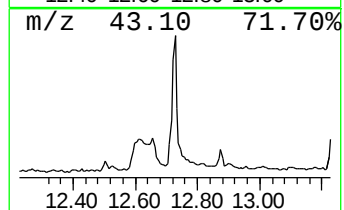
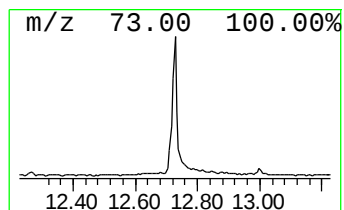
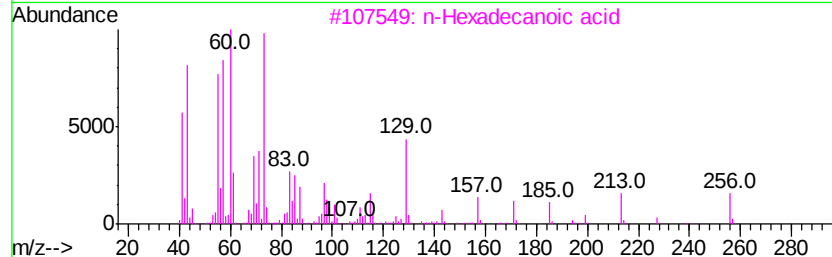
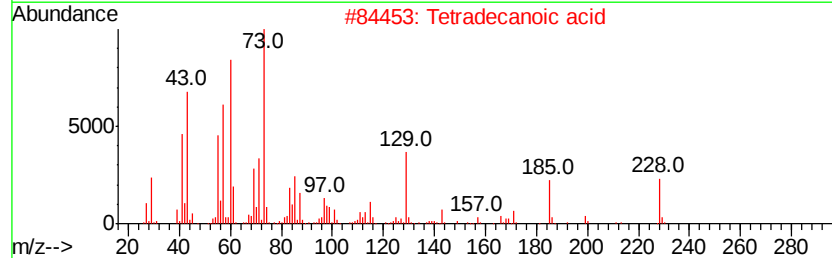
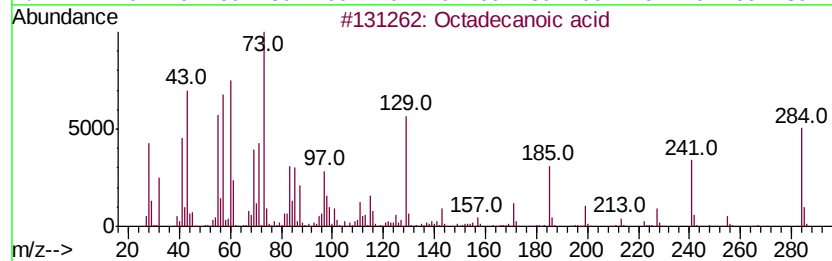
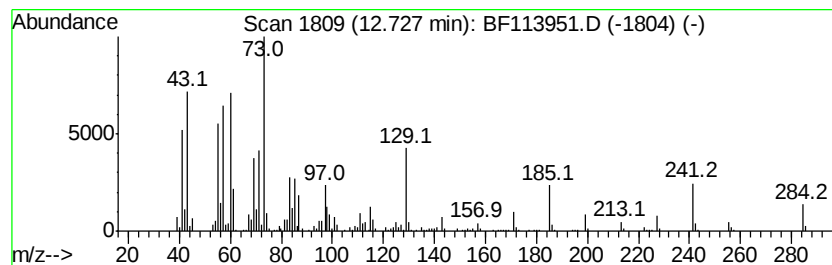
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.73	12.84 ng	786762	Phenanthrene-d10	11.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	74
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	58
4		n-Decanoic acid	172	C10H20O2	000334-48-5	53
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042419\
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Sample : K2522-07
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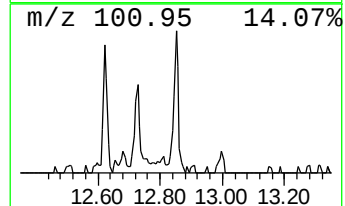
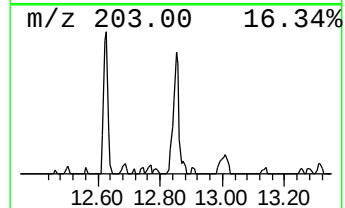
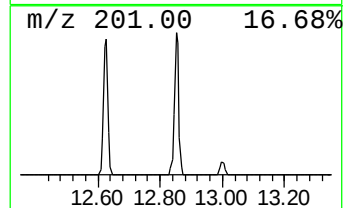
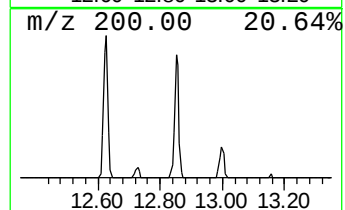
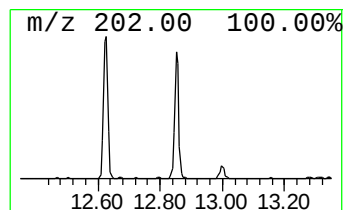
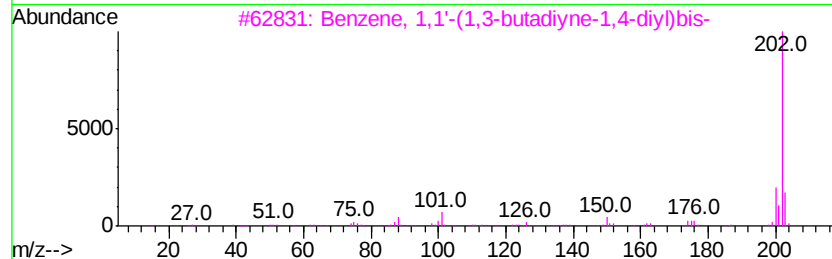
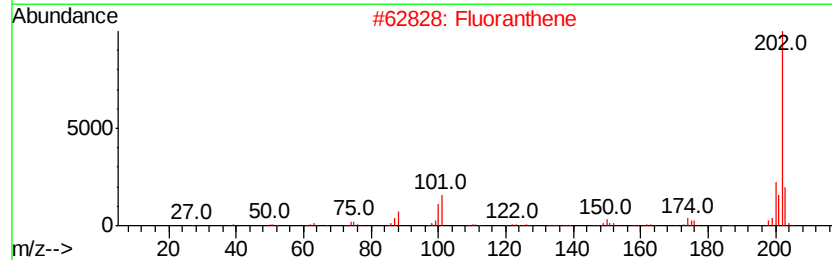
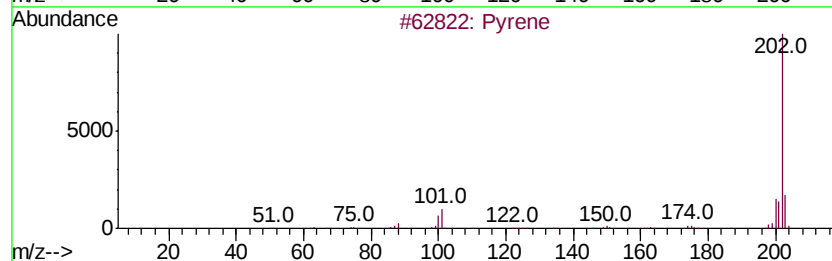
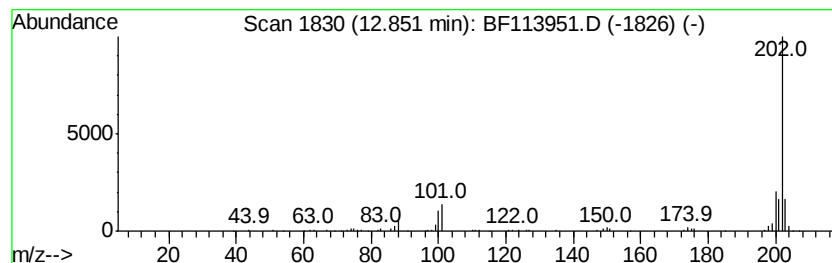
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzene, 1,1'-(1,3-butadiyn... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.85	2.00 ng	117638	Chrysene-d12	14.05
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Pyrene	202 C16H10	000129-00-0 94
2		Fluoranthene	202 C16H10	000206-44-0 91
3		Benzene, 1,1'-(1,3-butadiyne-1,4...	202 C16H10	000886-66-8 50
4		Furan-2-one, 2,5-dihydro-5-(4-me...	202 C12H10O3	024962-84-3 37
5		l-Leucyl-l-leucine, N,N'-dimethy...	528 C29H56N2O6	1000328-59-7 33



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042419\
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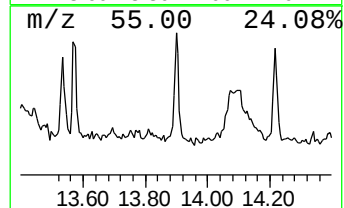
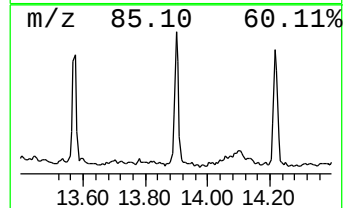
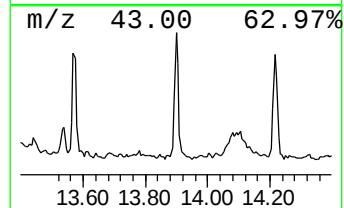
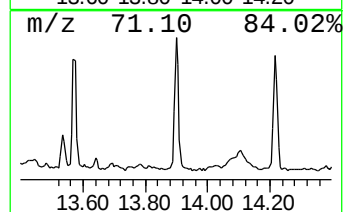
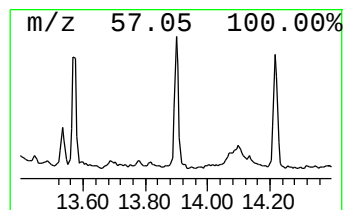
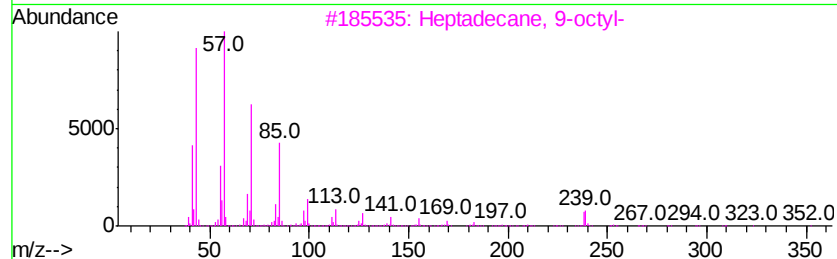
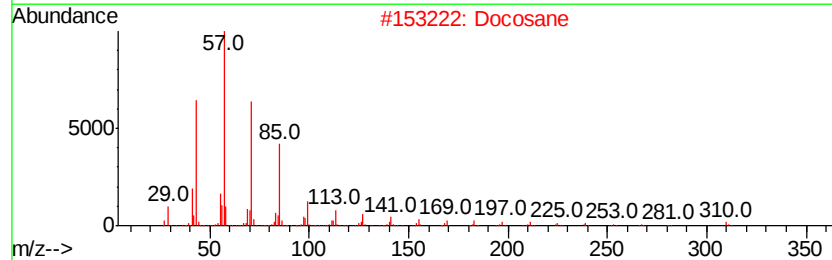
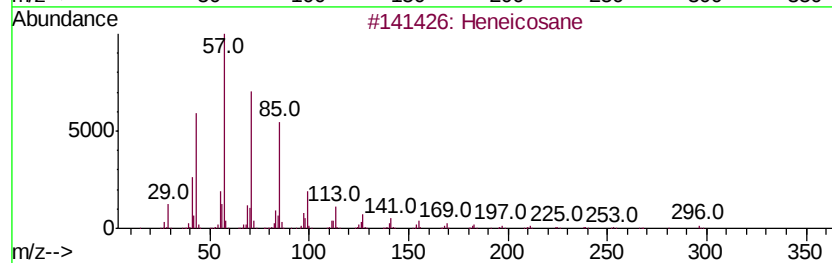
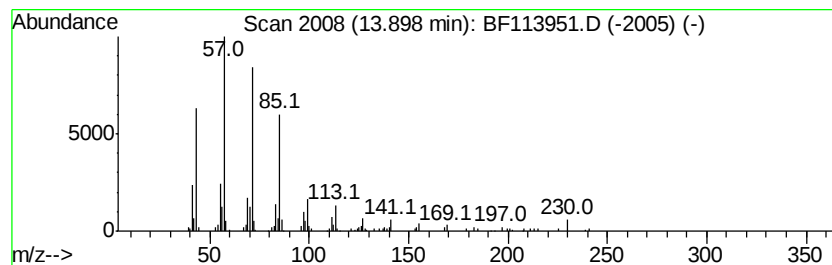
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heneicosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.90	2.83 ng	166414	Chrysene-d12	14.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Docosane	310	C22H46	000629-97-0	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		triacontane	422	C30H62	000638-68-6	91
5		Hexacosane	366	C26H54	000630-01-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042419\
Data File : BF113951.D
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Misc :
ALS Vial : 3 Sample Multiplier: 1

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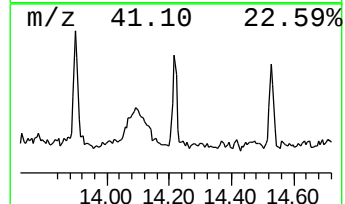
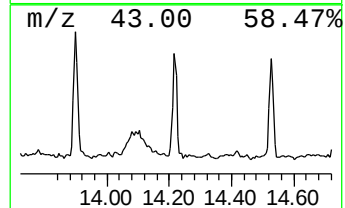
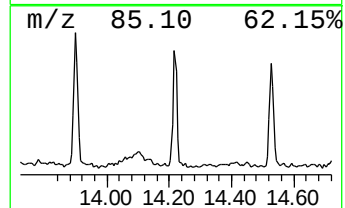
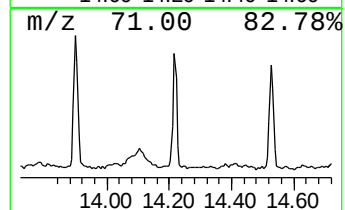
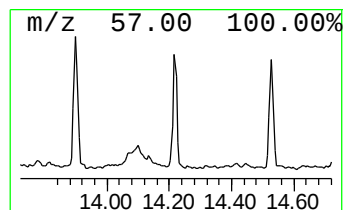
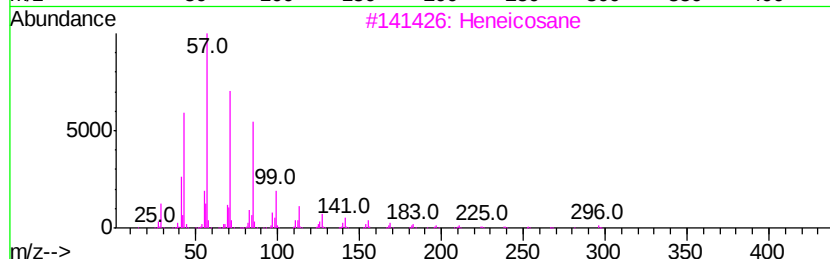
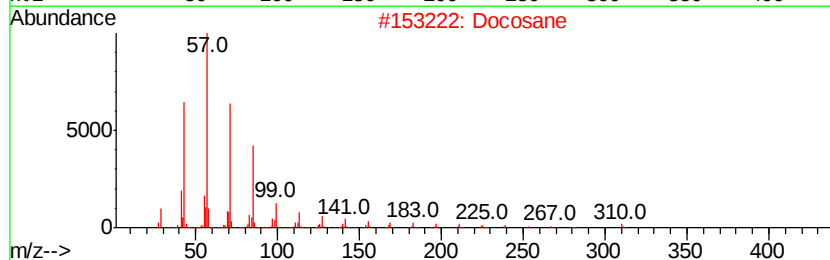
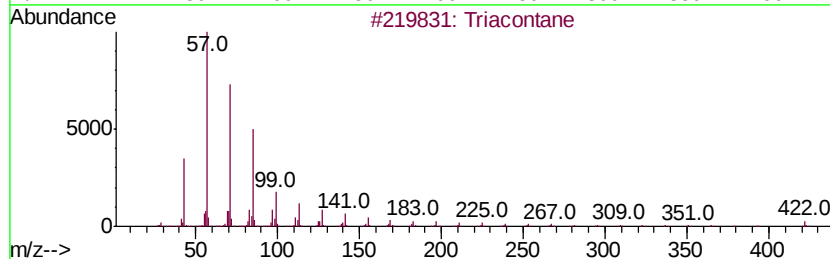
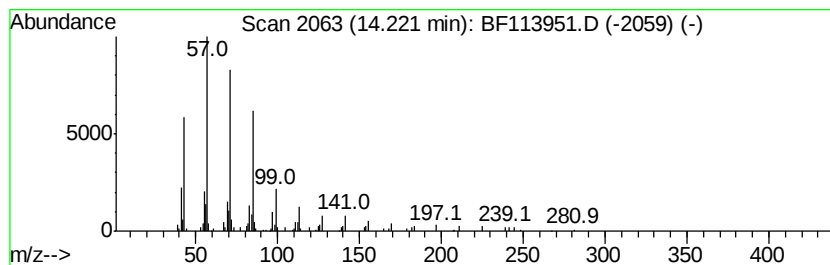
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Triacontane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.22	2.18 ng	128363	Chrysene-d12	14.05

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triacontane	422	C30H62	000638-68-6	91
2			Docosane	310	C22H46	000629-97-0	91
3			Heneicosane	296	C21H44	000629-94-7	91
4			Hexacosane	366	C26H54	000630-01-3	91
5			Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042419\
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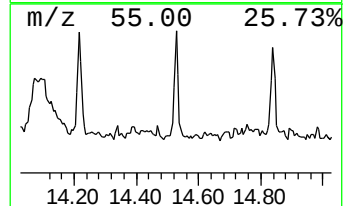
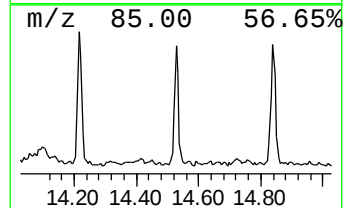
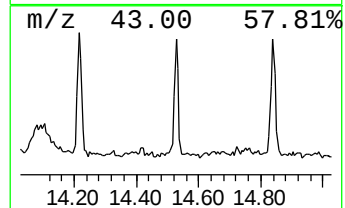
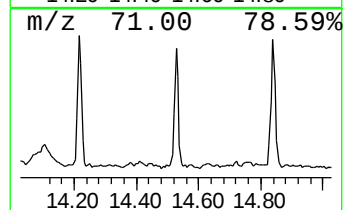
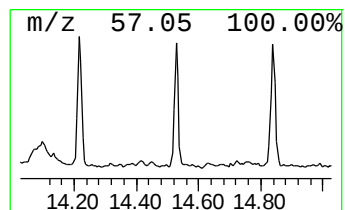
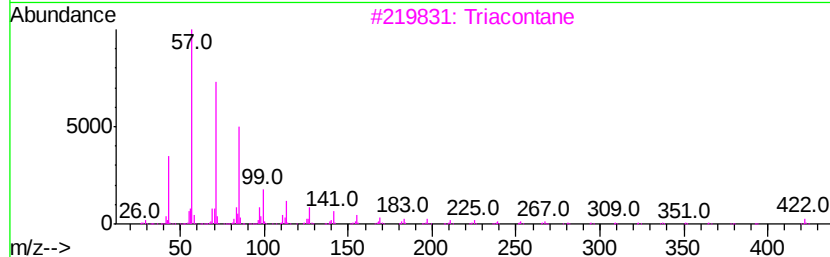
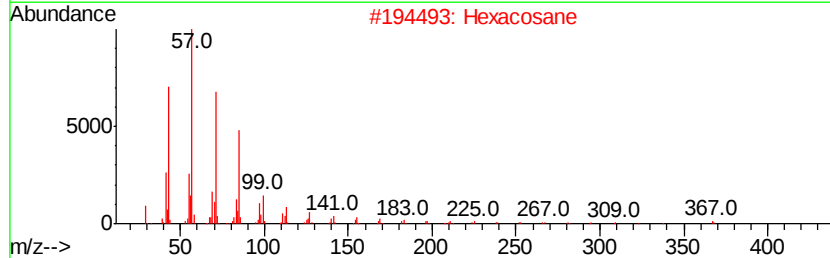
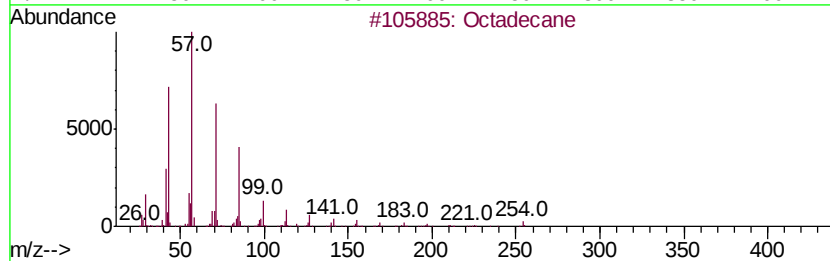
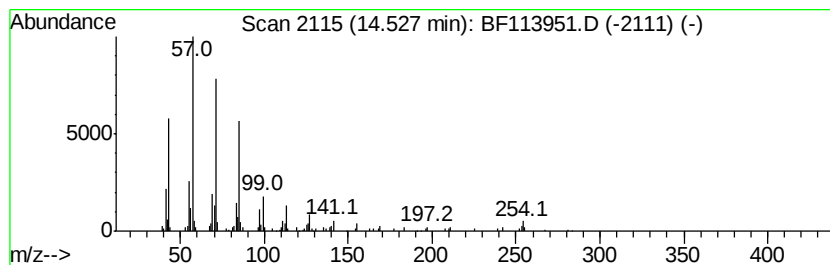
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Octadecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.53	2.28 ng	133879	Chrysene-d12	14.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	93
2		Hexacosane	366	C26H54	000630-01-3	90
3		triacontane	422	C30H62	000638-68-6	90
4		Heneicosane	296	C21H44	000629-94-7	90
5		2-methyloctacosane	408	C29H60	1000376-72-8	90



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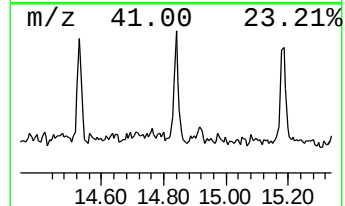
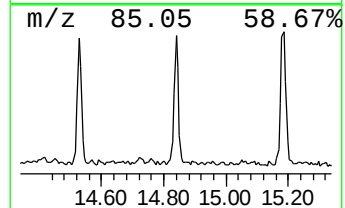
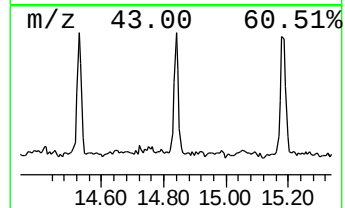
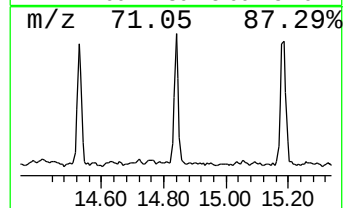
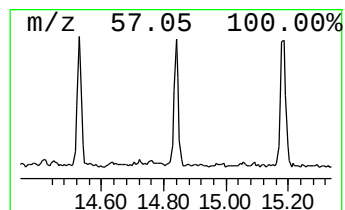
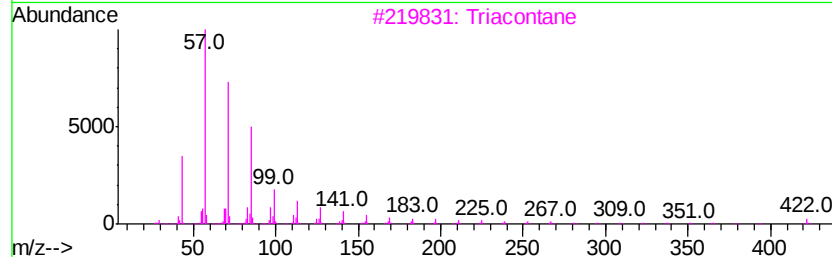
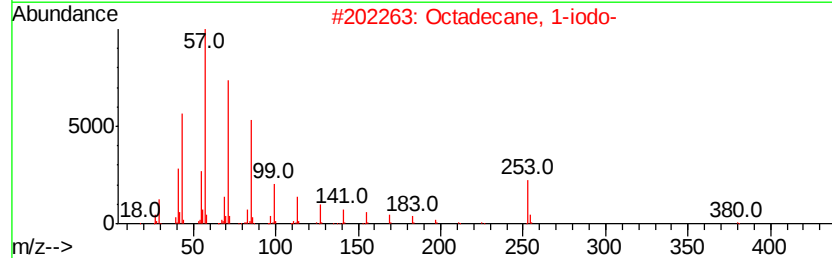
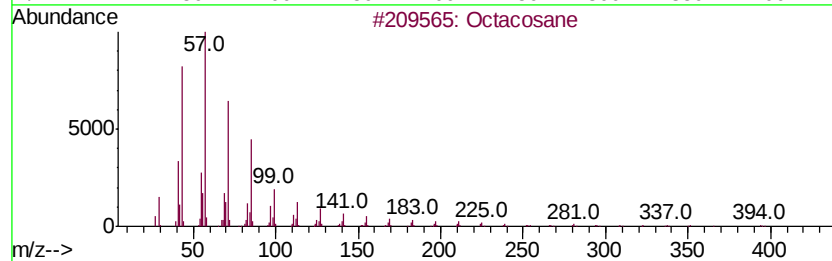
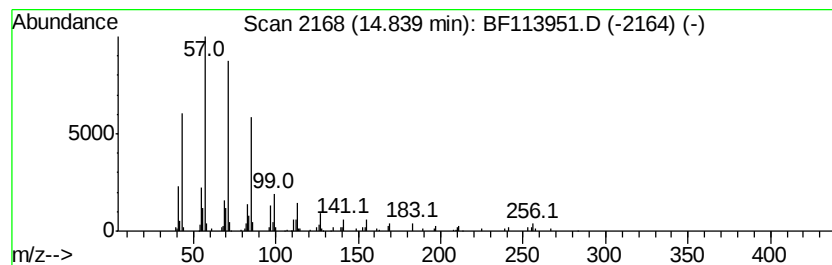
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Octadecane, 1-iodo- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.84	3.16 ng	142816	Perylene-d12	15.54

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	91
2			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	90
3			Triacontane	422	C30H62	000638-68-6	90
4			Heneicosane	296	C21H44	000629-94-7	90
5			Eicosane, 2-methyl-	296	C21H44	001560-84-5	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042419\
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ClientSampleId :
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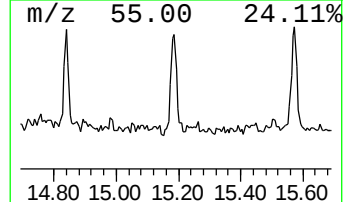
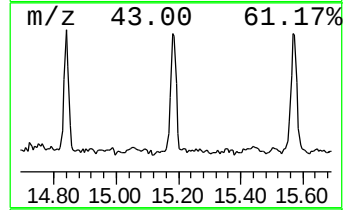
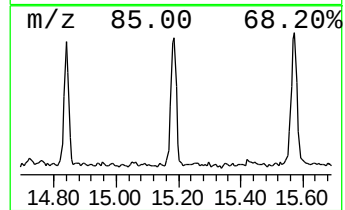
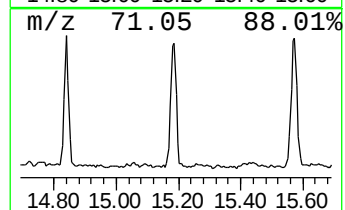
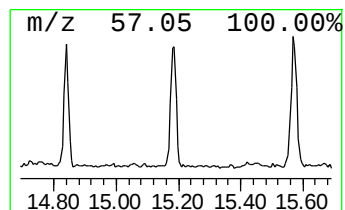
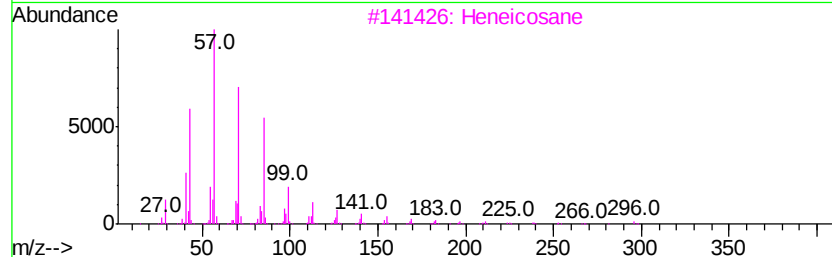
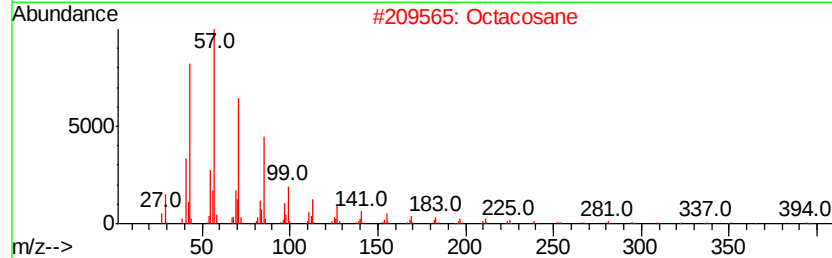
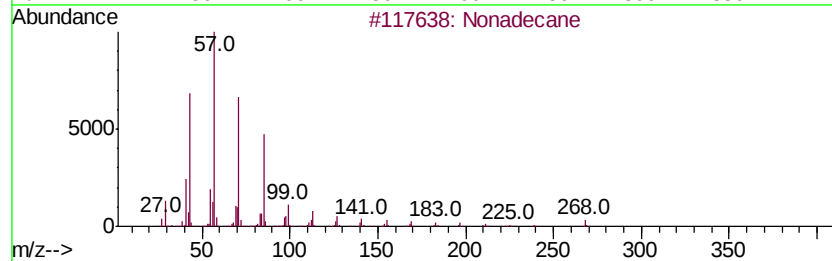
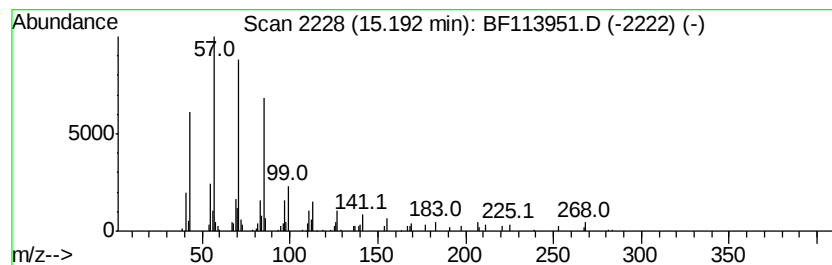
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Nonadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.19	3.65 ng	164701	Perylene-d12	15.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	96
2		Octacosane	394	C28H58	000630-02-4	91
3		Heneicosane	296	C21H44	000629-94-7	90
4		Triacontane	422	C30H62	000638-68-6	90
5		2-methyloctacosane	408	C29H60	1000376-72-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042419\
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 Acq On : 24 Apr 2019 10:46
 Operator : JU/SJ
 Sample : K2522-07
 Misc :
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Instrument :
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 ClientSampleId :
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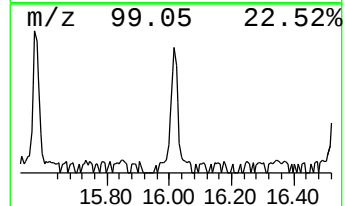
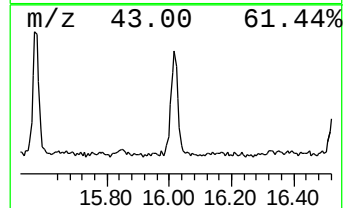
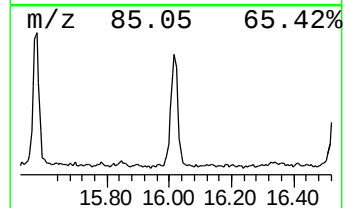
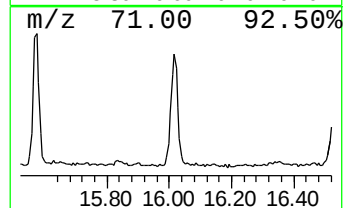
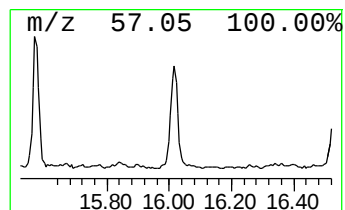
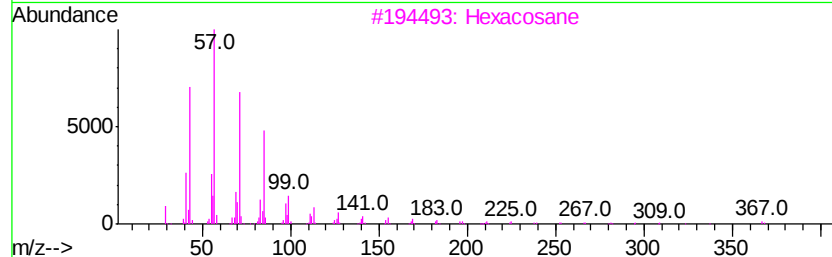
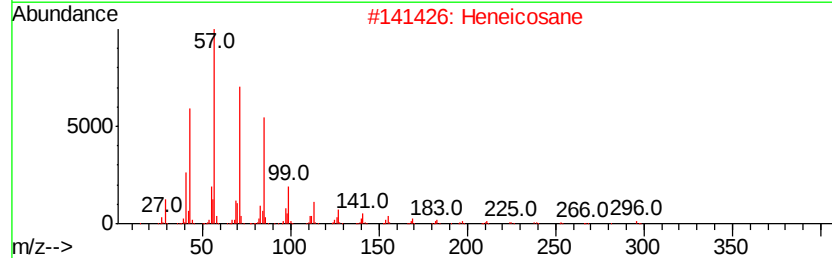
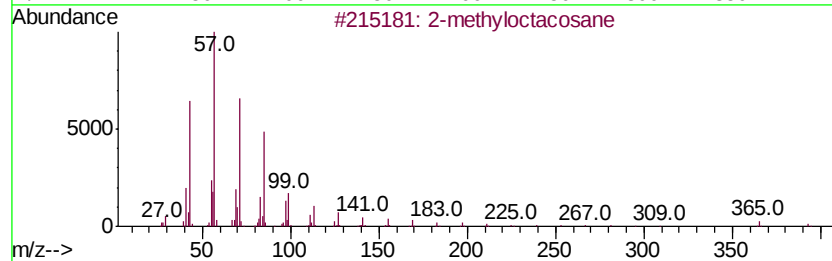
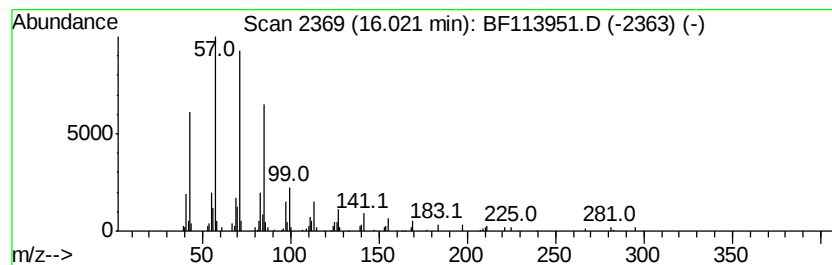
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 15 2-methyloctacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.02	3.74 ng	168786	Perylene-d12	15.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-methyloctacosane	408	C29H60	1000376-72-8	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Hexacosane	366	C26H54	000630-01-3	91
4		triacontane	422	C30H62	000638-68-6	91
5		Heptadecane	240	C17H36	000629-78-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042419\
Data File : BF113951.D
Acq On : 24 Apr 2019 10:46
Operator : JU/SJ
Sample : K2522-07
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP-WOODBURY

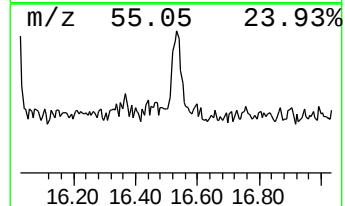
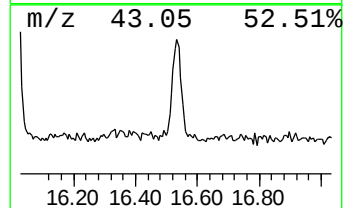
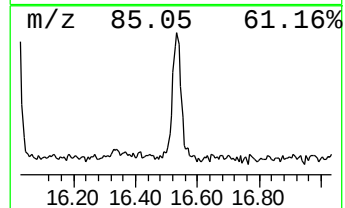
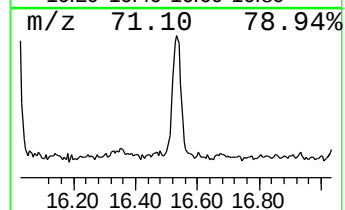
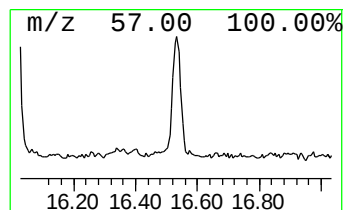
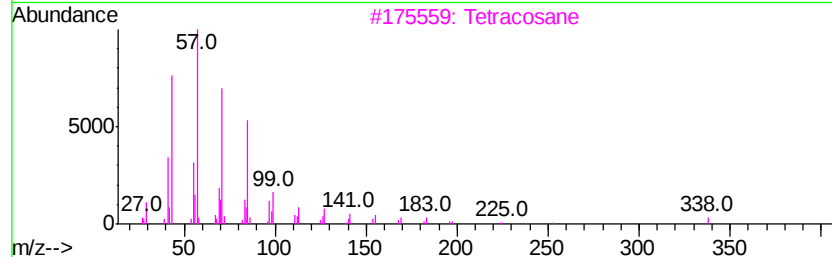
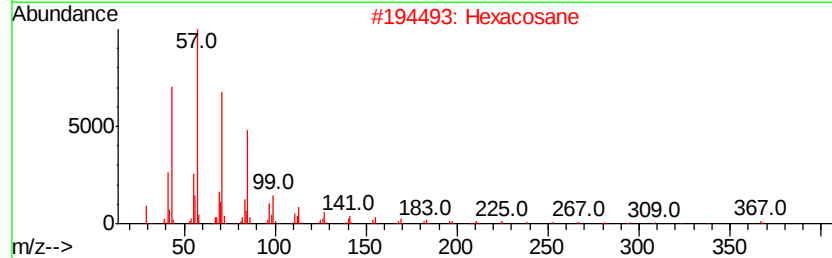
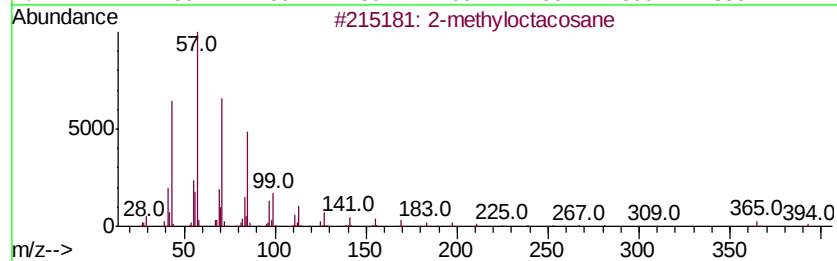
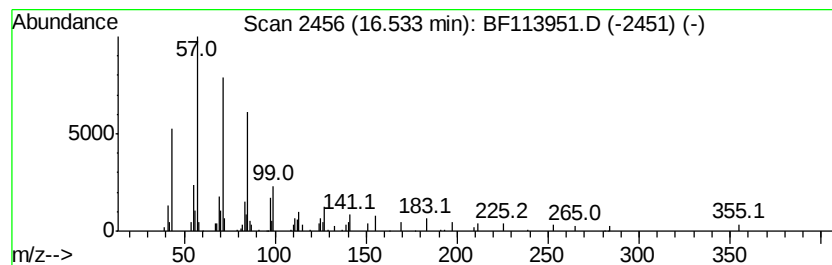
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Eicosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.53	2.44 ng	110336	Perylene-d12	15.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-methyloctacosane	408	C29H60	1000376-72-8	91
2		Hexacosane	366	C26H54	000630-01-3	90
3		Tetracosane	338	C24H50	000646-31-1	87
4		Eicosane	282	C20H42	000112-95-8	87
5		triacontane	422	C30H62	000638-68-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042419\
Data File : BF113951.D
Acq On : 24 Apr 2019 10:46
Operator : JU/SJ
Sample : K2522-07
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP-WOODBURY

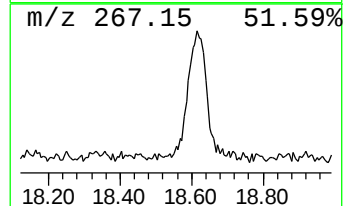
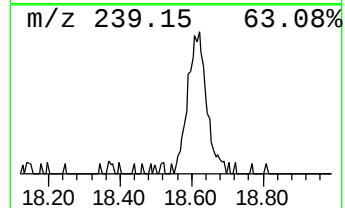
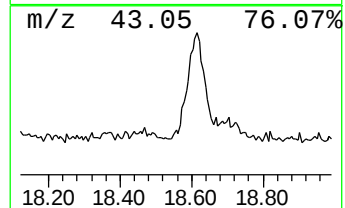
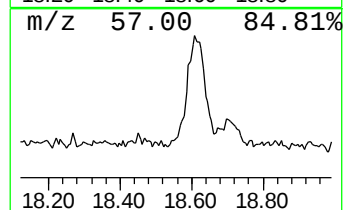
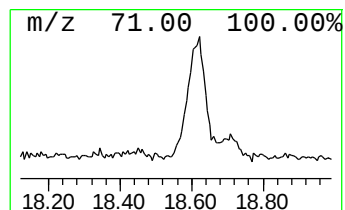
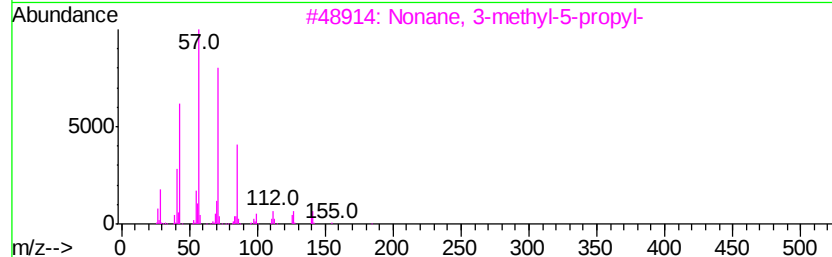
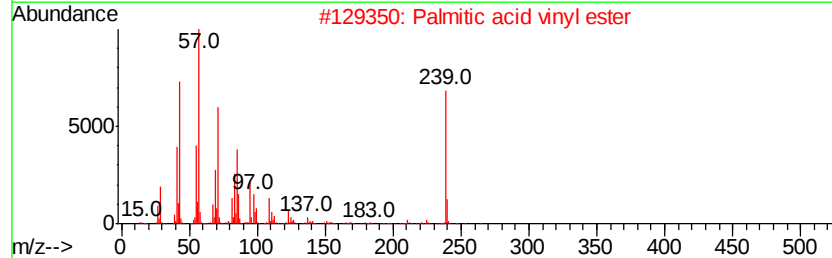
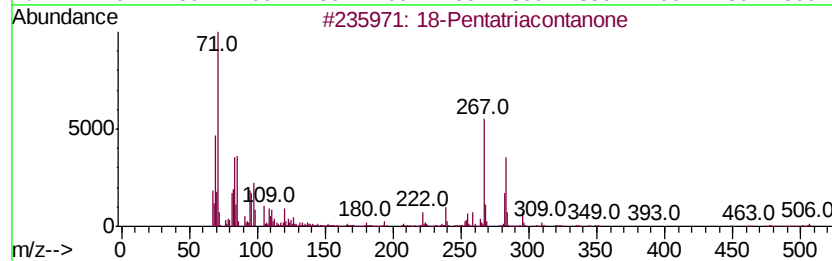
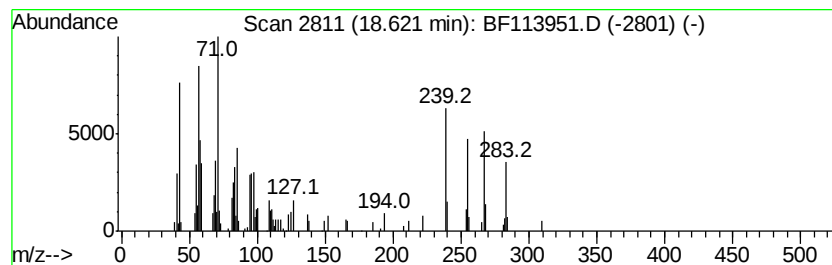
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 18-Pentatriacontanone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.62	5.72 ng	258378	Perylene-d12	15.54

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			18-Pentatriacontanone	507	C35H70O	000504-53-0	58
2			Palmitic acid vinyl ester	282	C18H34O2	000693-38-9	30
3			Nonane, 3-methyl-5-propyl-	184	C13H28	031081-18-2	27
4			Nonane, 1-iodo-	254	C9H19I	004282-42-2	27
5			Nonane, 2-methyl-5-propyl-	184	C13H28	031081-17-1	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF042419\
 Data File : BF113951.D
 Acq On : 24 Apr 2019 10:46
 Operator : JU/SJ
 Sample : K2522-07
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-WOODBURY

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF042219.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...	2.21	6.3	ng	231984	1	6.90	739910	20.0
unknown6.67	6.67	86.3	ng	3191410	1	6.90	739910	20.0
2,4,7,9-Tetrameth...	9.37	2.9	ng	166070	3	9.93	1129020	20.0
n-Hexadecanoic acid	11.95	15.8	ng	966832	4	11.42	1225310	20.0
2-Ethylbutyric ac...	12.66	3.4	ng	207090	4	11.42	1225310	20.0
Octadecanoic acid	12.73	12.8	ng	786762	4	11.42	1225310	20.0
Benzene, 1,1'-(1,...	12.85	2.0	ng	117638	5	14.05	1175550	20.0
Heneicosane	13.90	2.8	ng	166414	5	14.05	1175550	20.0
triacontane	14.22	2.2	ng	128363	5	14.05	1175550	20.0
Octadecane	14.53	2.3	ng	133879	5	14.05	1175550	20.0
Octadecane, 1-iodo-	14.84	3.2	ng	142816	6	15.54	903525	20.0
Nonadecane	15.19	3.6	ng	164701	6	15.54	903525	20.0
2-methyloctacosane	16.02	3.7	ng	168786	6	15.54	903525	20.0
Eicosane	16.53	2.4	ng	110336	6	15.54	903525	20.0
18-Pentatriaconta...	18.62	5.7	ng	258378	6	15.54	903525	20.0