

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF050219\
 Data File : BF114081.D
 Acq On : 2 May 2019 13:02
 Operator : JU/SJ
 Sample : K2584-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSSI-0419-S-DH-7

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	3.975	315	321	330	rVB	89199	123312	2.09%	0.367%
2	5.516	578	583	586	rBV	2896969	3039517	51.44%	9.052%
3	6.528	749	755	758	rBV	2874057	3196400	54.10%	9.519%
4	6.663	773	778	781	rBV	3370511	3321068	56.21%	9.891%
5	6.887	813	816	820	rBV	553152	436888	7.39%	1.301%
6	7.051	839	844	847	rBV	2642998	2540655	43.00%	7.566%
7	7.451	906	912	915	rBV	2095904	2008145	33.99%	5.981%
8	8.169	1030	1034	1041	rBV	487071	409427	6.93%	1.219%
9	9.245	1213	1217	1220	rBV	3521097	3367687	57.00%	10.029%
10	9.634	1279	1283	1288	rBV	327407	255760	4.33%	0.762%
11	9.922	1329	1332	1337	rBV	570236	536960	9.09%	1.599%
12	10.716	1462	1467	1470	rBV	2572110	2650958	44.87%	7.895%
13	11.410	1581	1585	1588	rBV	729636	614080	10.39%	1.829%
14	11.933	1668	1674	1682	rBV	316689	389740	6.60%	1.161%
15	12.622	1788	1791	1800	rBV	156690	149513	2.53%	0.445%
16	12.722	1803	1808	1815	rBV	584063	639892	10.83%	1.906%
17	12.851	1827	1830	1834	rVB	130651	123909	2.10%	0.369%
18	12.998	1850	1855	1858	rBV	5027183	5908493	100.00%	17.596%
19	13.892	2004	2007	2022	rVB	272769	416400	7.05%	1.240%
20	14.057	2028	2035	2038	rBV	668912	707848	11.98%	2.108%
21	14.216	2059	2062	2067	rBV	85052	90234	1.53%	0.269%
22	14.533	2111	2116	2127	rBV	189614	278433	4.71%	0.829%
23	14.845	2165	2169	2174	rBV	247750	273933	4.64%	0.816%
24	15.121	2212	2216	2219	rBV	40269	65609	1.11%	0.195%
25	15.192	2224	2228	2232	rBV	332599	369864	6.26%	1.102%
26	15.557	2285	2290	2292	rBV	451378	599591	10.15%	1.786%
27	15.580	2292	2294	2300	rVB	286777	322879	5.46%	0.962%
28	16.027	2364	2370	2376	rBV	205475	317297	5.37%	0.945%
29	16.098	2376	2382	2395	rVB	35661	123543	2.09%	0.368%
30	16.539	2452	2457	2462	rBV	80662	139998	2.37%	0.417%
31	17.145	2556	2560	2568	rVB2	35476	68291	1.16%	0.203%
32	17.680	2644	2651	2657	rBV2	26452	91643	1.55%	0.273%

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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-BF042219.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

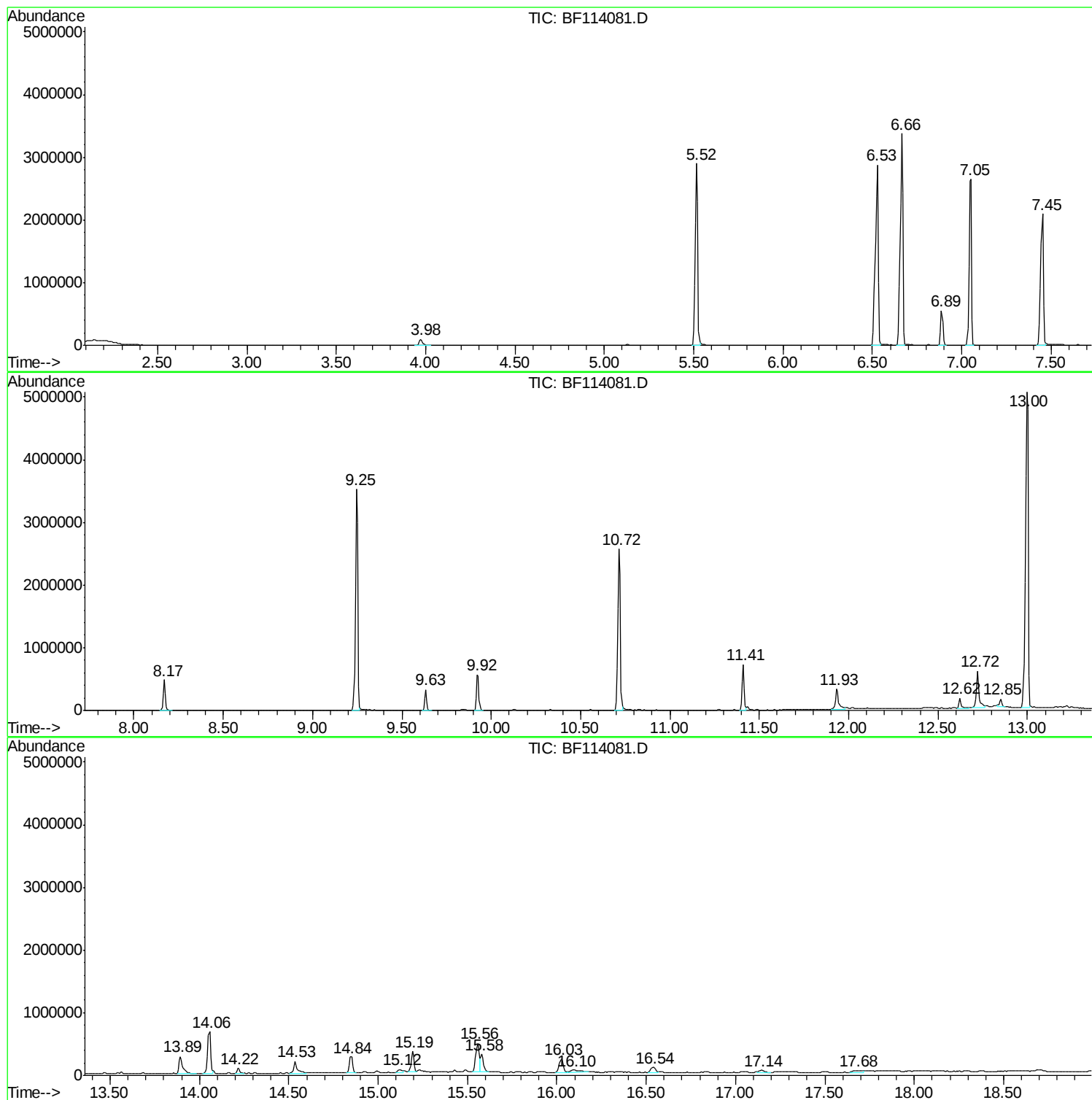
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF042219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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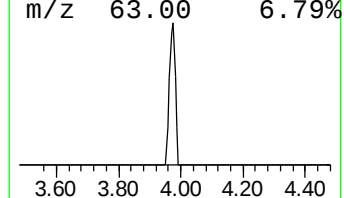
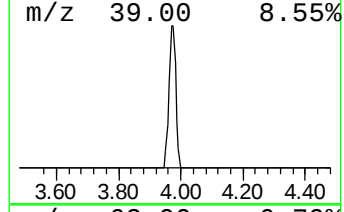
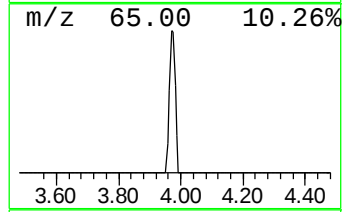
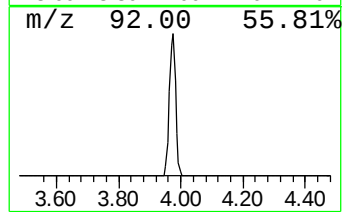
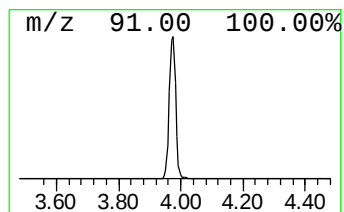
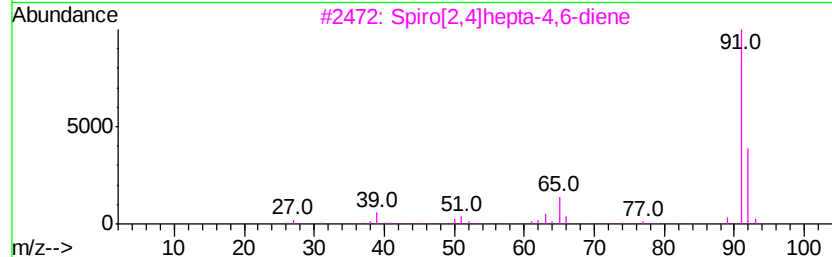
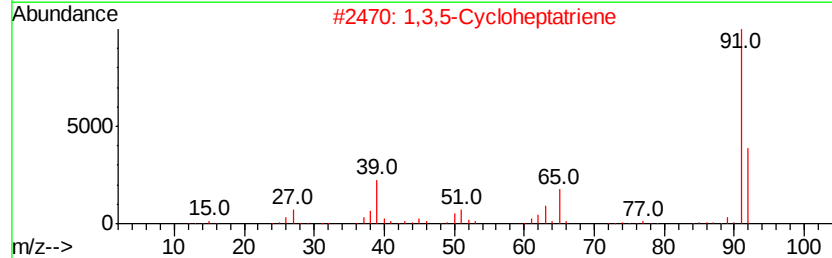
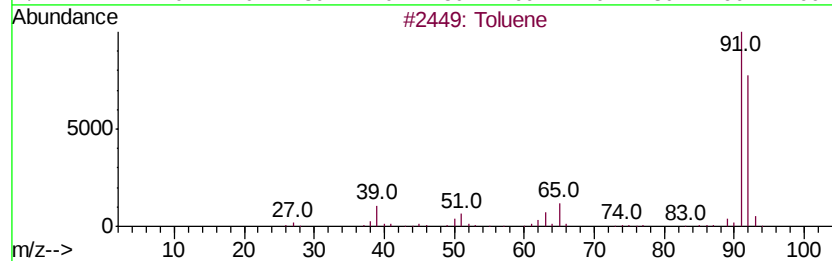
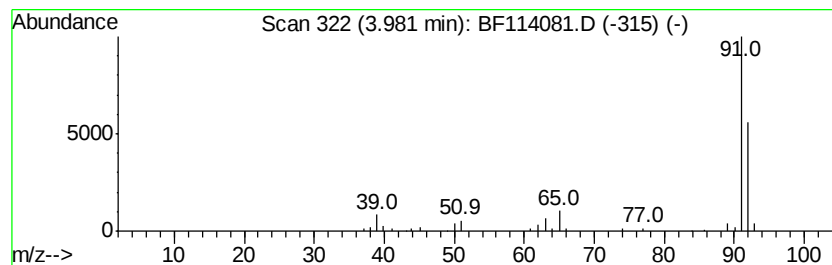
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 1,3,5-Cycloheptatriene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.98	5.65 ng	123312	1,4-Dichlorobenzene-d4	6.89

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Toluene	92	C7H8	000108-88-3	95
2			1,3,5-Cycloheptatriene	92	C7H8	000544-25-2	90
3			Spiro[2.4]hepta-4,6-diene	92	C7H8	000765-46-8	68
4			Molybdenum, di-.mu.-chlorobis[(1...	532	C20H26Cl2Mo2	035625-66-2	64
5			2,5-Norbornadiene	92	C7H8	000121-46-0	64



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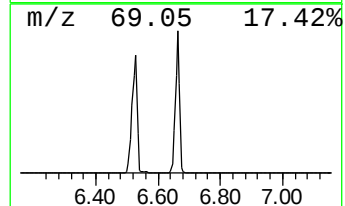
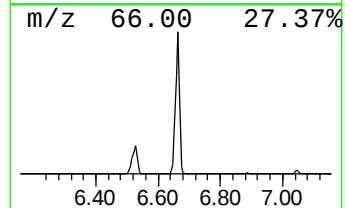
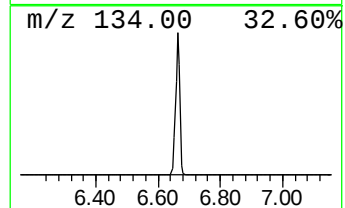
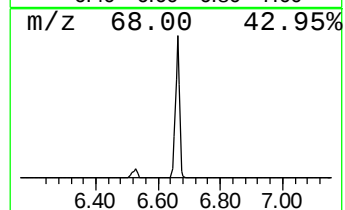
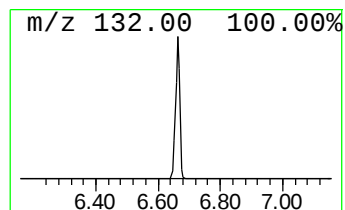
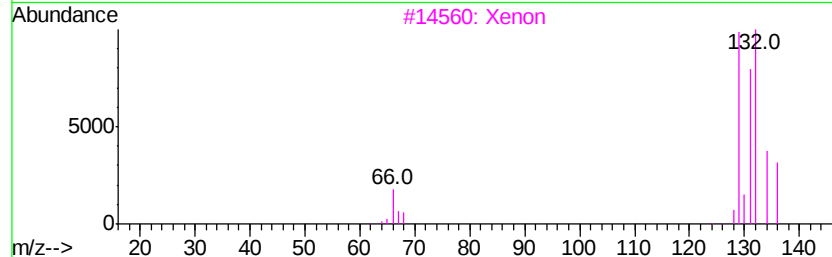
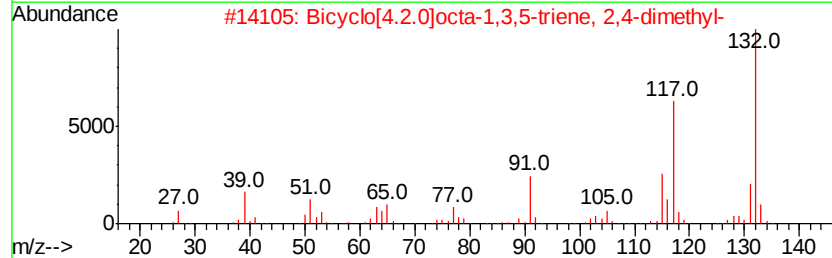
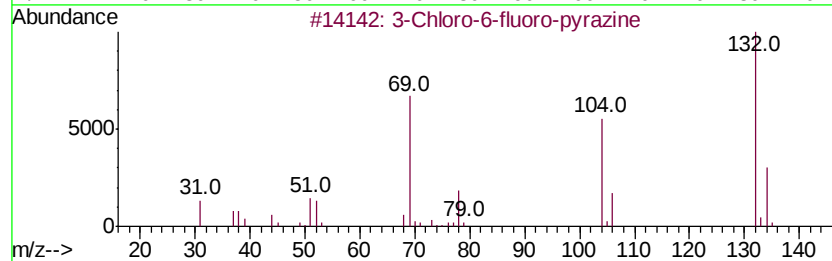
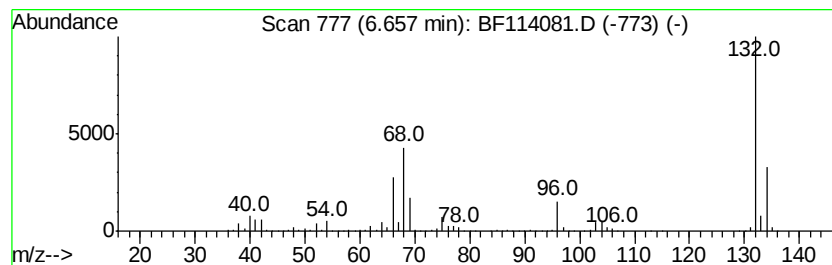
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.66 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.66	152.03 ng	3321070	1,4-Dichlorobenzene-d4	6.89

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine			132	C4H2ClFN2	1000146-10-7	35
2	Bicyclo[4.2.0]octa-1,3,5-triene, ...			132	C10H12	028749-81-7	9
3	Xenon			132	Xe	007440-63-3	9
4	4-Chloro-2-fluoro-pyrimidine			132	C4H2ClFN2	051422-00-5	9
5	5-Fluoro-2-chloropyrimidine			132	C4H2ClFN2	062802-42-0	9



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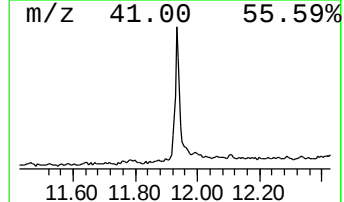
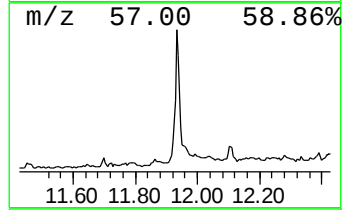
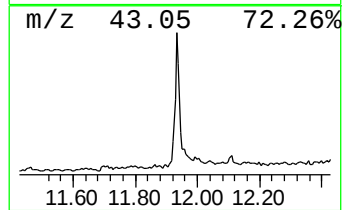
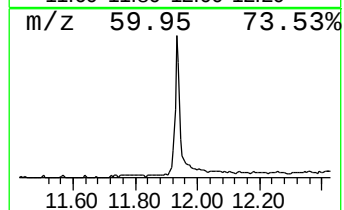
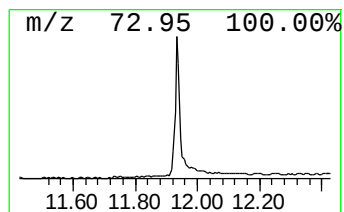
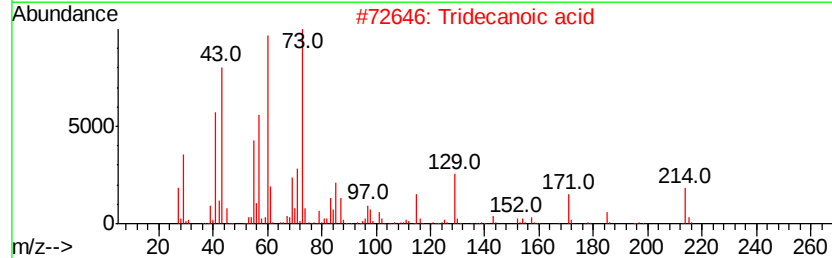
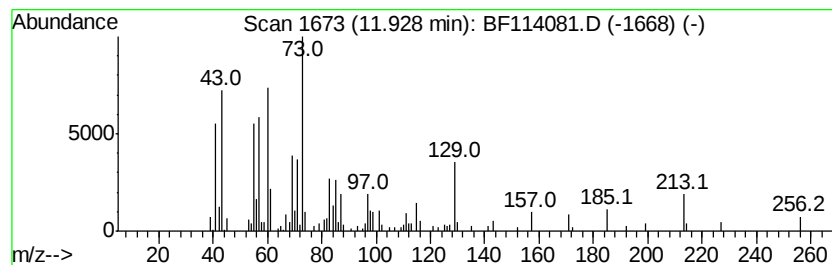
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF042219.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	12.69 ng	389740	Phenanthrene-d10	11.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Tridecanoic acid	214	C13H26O2	000638-53-9	89
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	76
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	52
5		n-Decanoic acid	172	C10H20O2	000334-48-5	35



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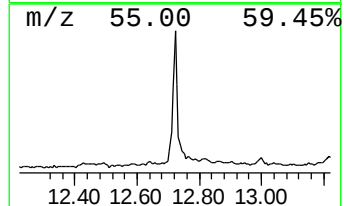
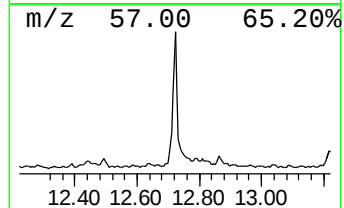
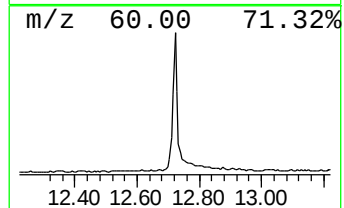
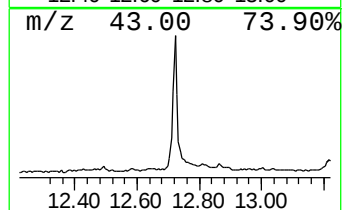
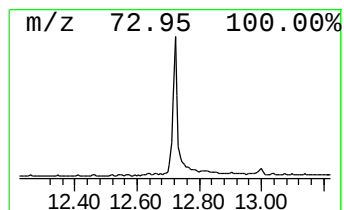
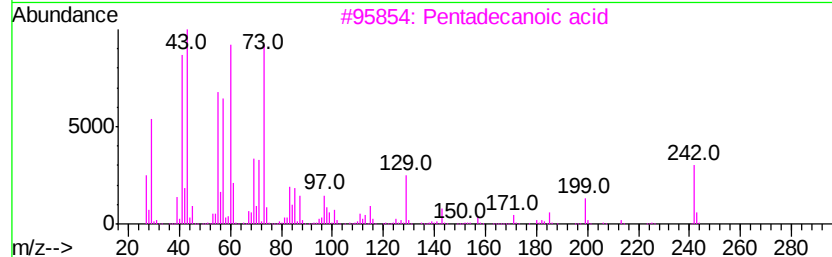
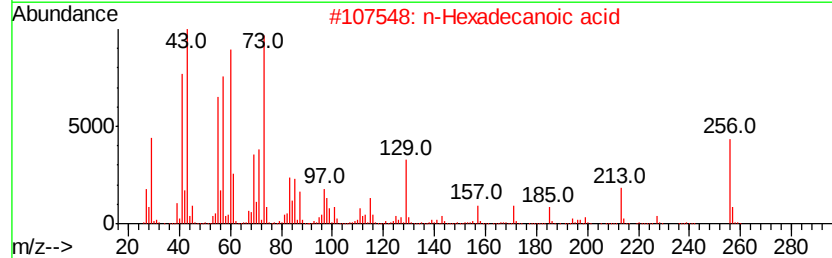
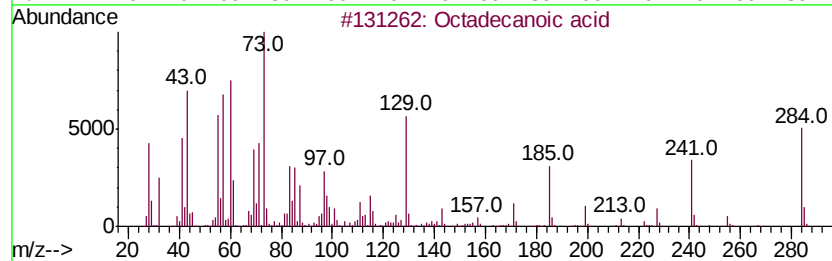
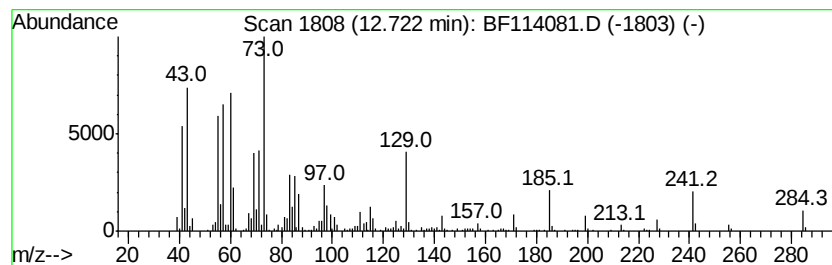
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Octadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.72	20.84 ng	639892	Phenanthrene-d10	11.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	83
5	n-Decanoic acid	172	C10H20O2	000334-48-5	58



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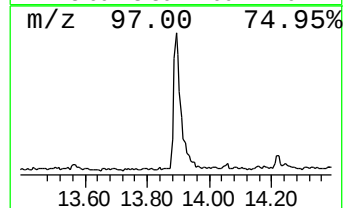
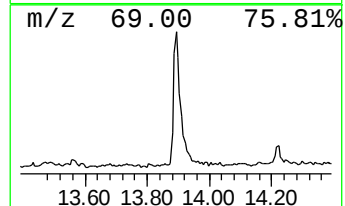
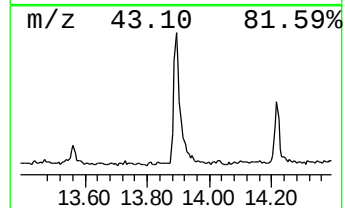
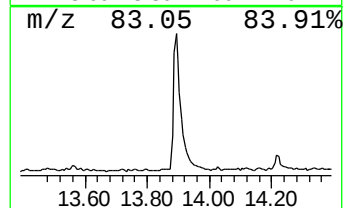
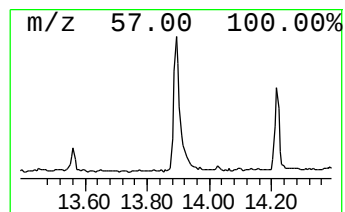
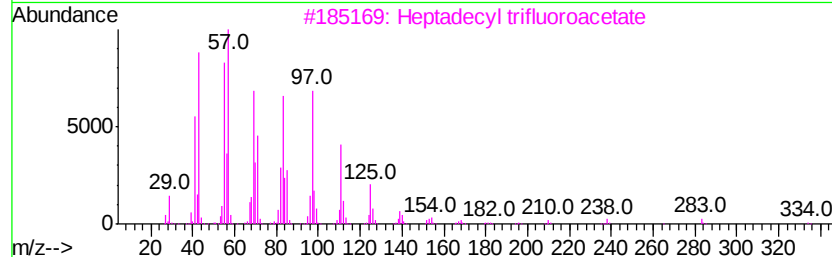
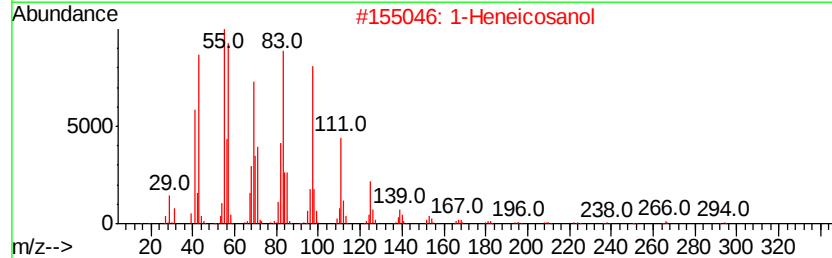
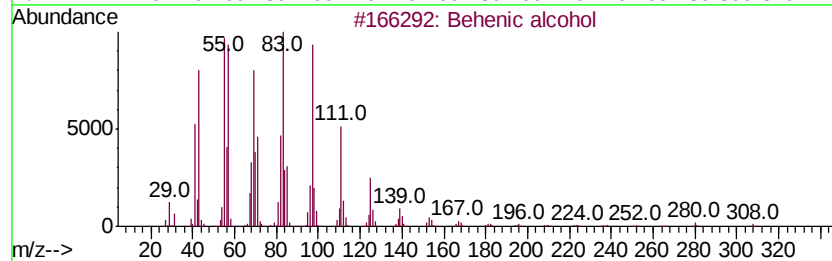
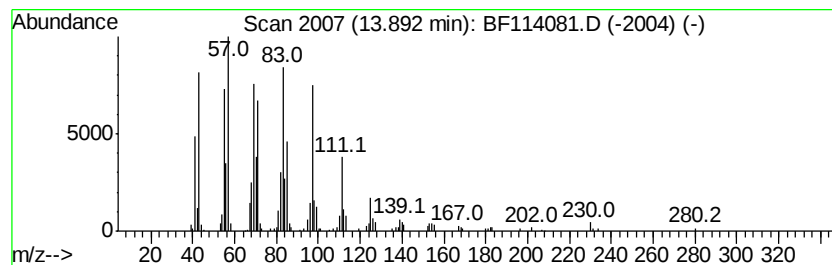
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Behenic alcohol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.89	11.77 ng	416400	Chrysene-d12	14.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Behenic alcohol	326	C22H46O	000661-19-8	93
2	1-Heneicosanol	312	C21H44O	015594-90-8	93
3	Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	93
4	Heptadecyl heptafluorobutyrte	452	C21H35F7O2	959085-66-6	93
5	Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	93



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SSSI-0419-S-DH-7

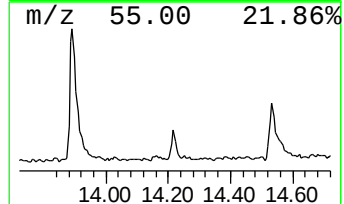
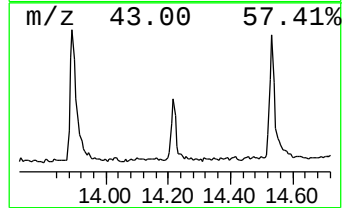
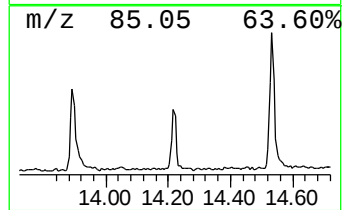
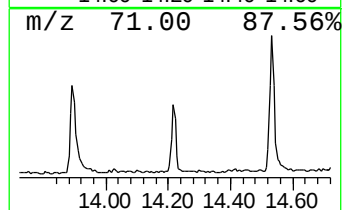
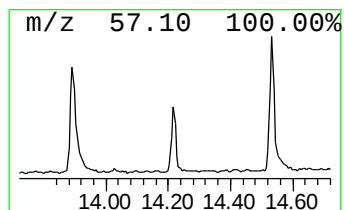
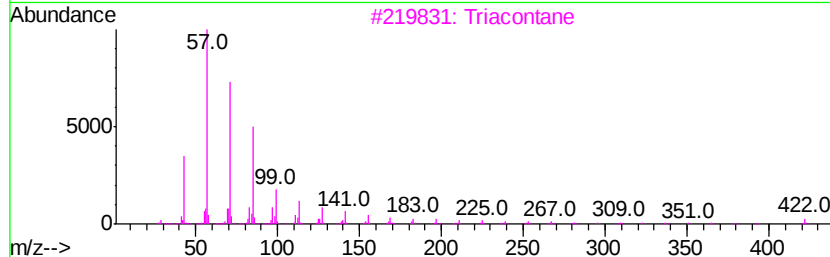
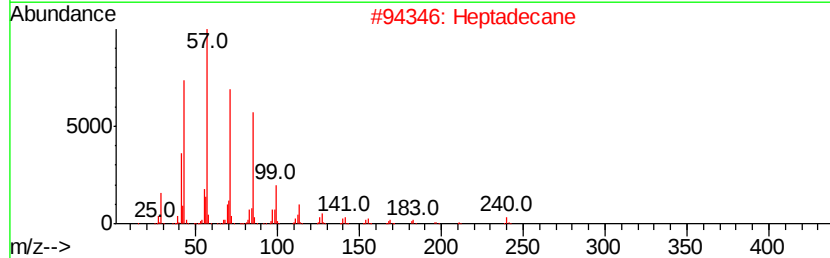
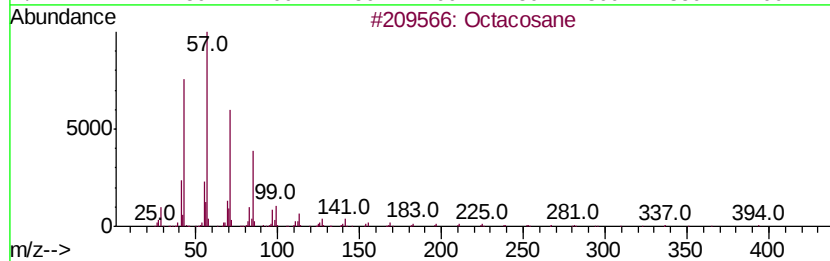
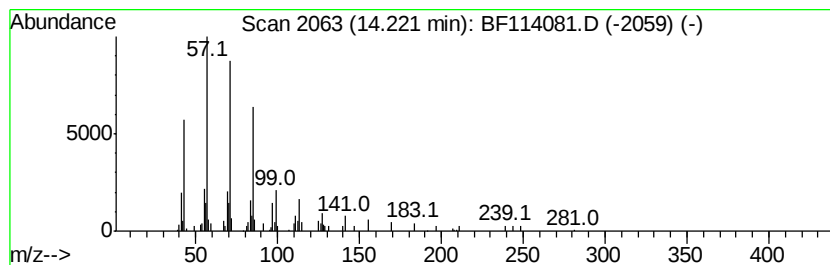
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octacosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.22	2.55 ng	90234	Chrysene-d12	14.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Heptadecane	240	C17H36	000629-78-7	91
3		Triacontane	422	C30H62	000638-68-6	91
4		Tetracosane	338	C24H50	000646-31-1	91
5		Docosane	310	C22H46	000629-97-0	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF050219\
Data File : BF114081.D
Acq On : 2 May 2019 13:02
Operator : JU/SJ
Sample : K2584-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SSSI-0419-S-DH-7

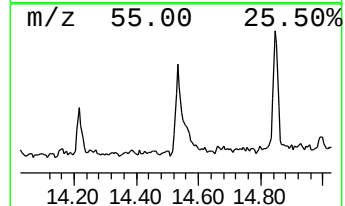
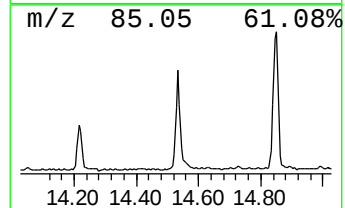
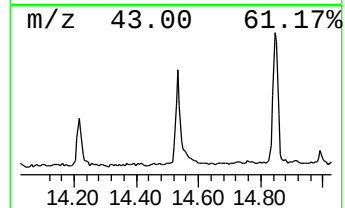
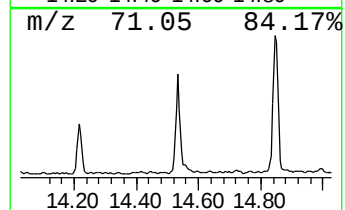
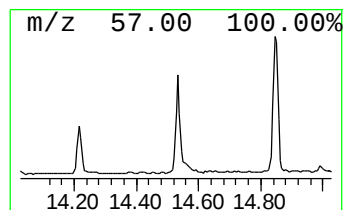
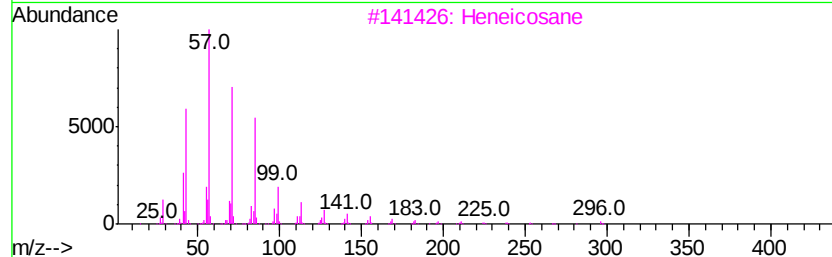
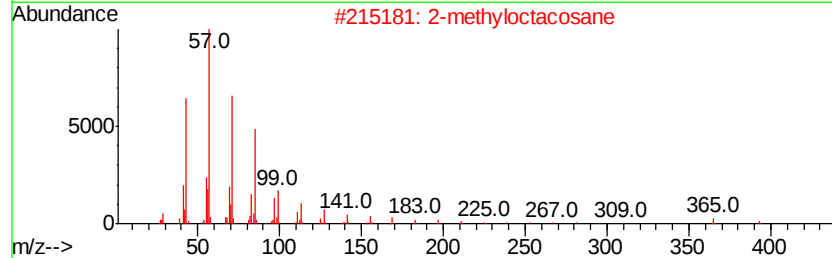
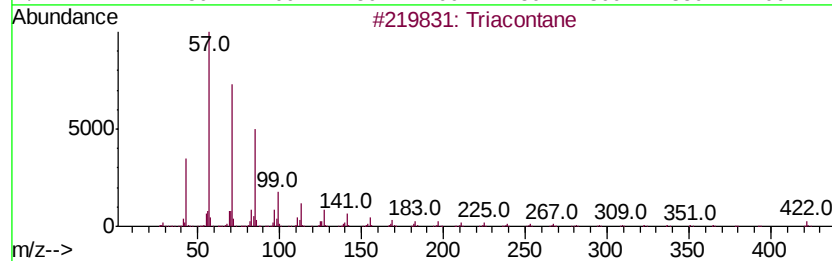
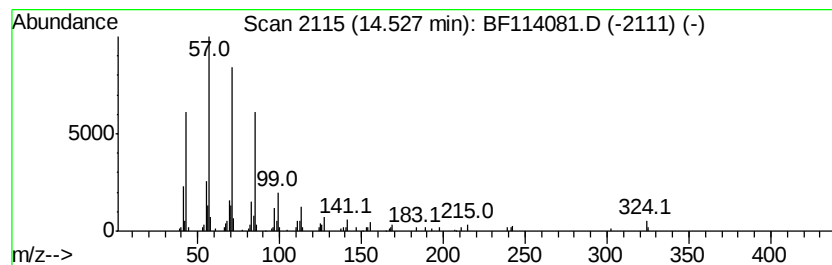
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Triacontane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.53	7.87 ng	278433	Chrysene-d12	14.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Triacontane	422	C30H62	000638-68-6	91
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Heneicosane	296	C21H44	000629-94-7	90
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
5		Hexacosane	366	C26H54	000630-01-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF050219\
Data File : BF114081.D
Acq On : 2 May 2019 13:02
Operator : JU/SJ
Sample : K2584-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SSSI-0419-S-DH-7

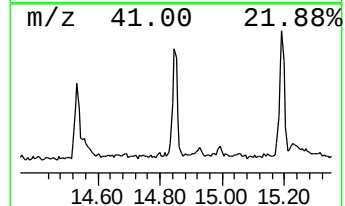
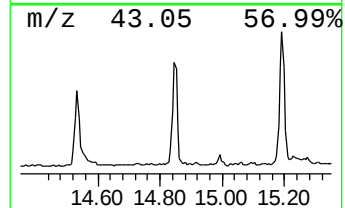
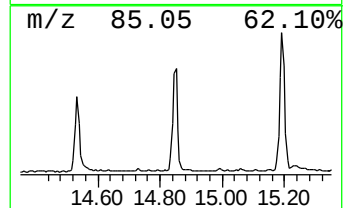
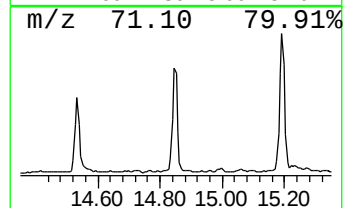
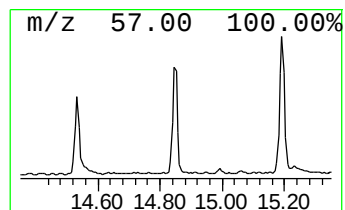
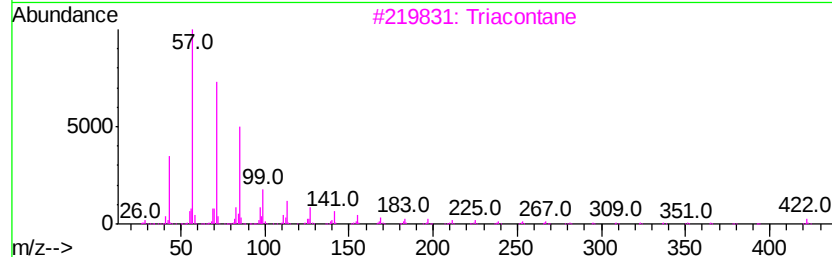
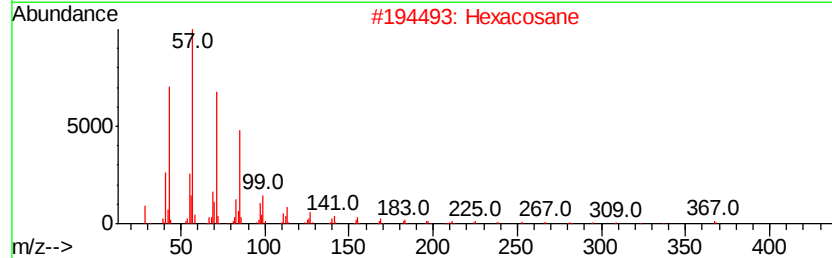
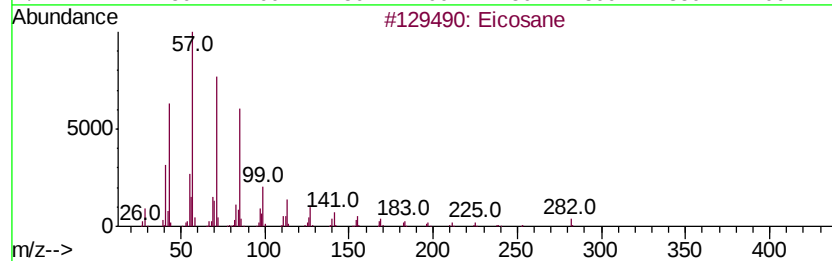
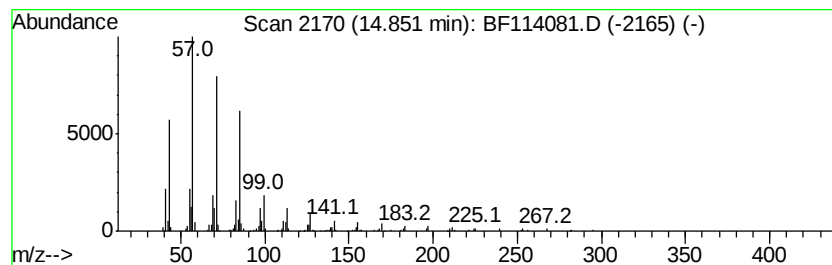
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Eicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	9.14 ng	273933	Perylene-d12	15.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	97
2	Hexacosane		366	C26H54	000630-01-3	93
3	Triacontane		422	C30H62	000638-68-6	91
4	2-methyloctacosane		408	C29H60	1000376-72-8	91
5	Heneicosane		296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF050219\
Data File : BF114081.D
Acq On : 2 May 2019 13:02
Operator : JU/SJ
Sample : K2584-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SSSI-0419-S-DH-7

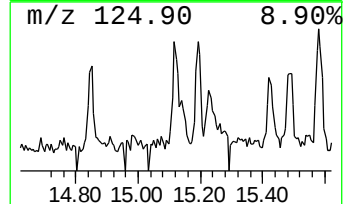
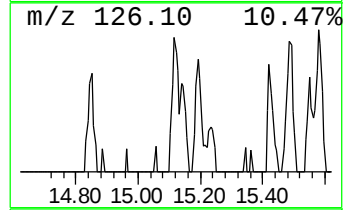
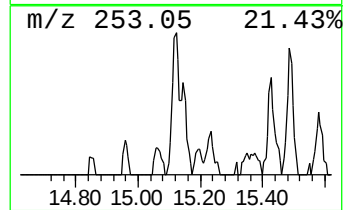
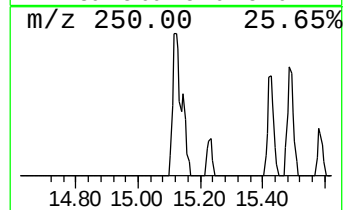
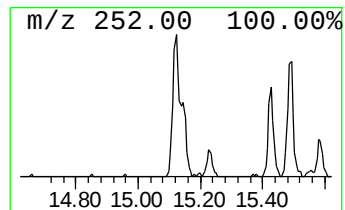
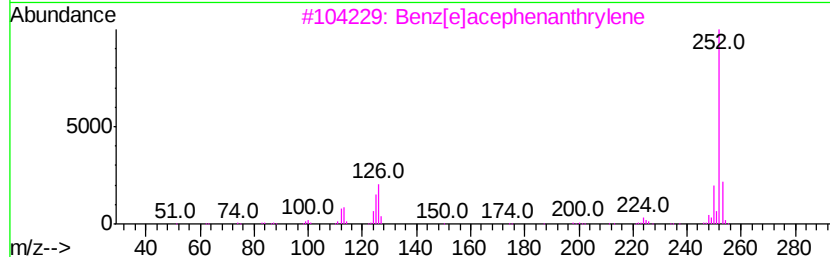
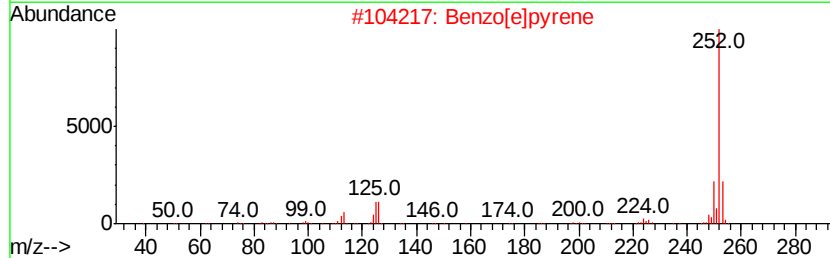
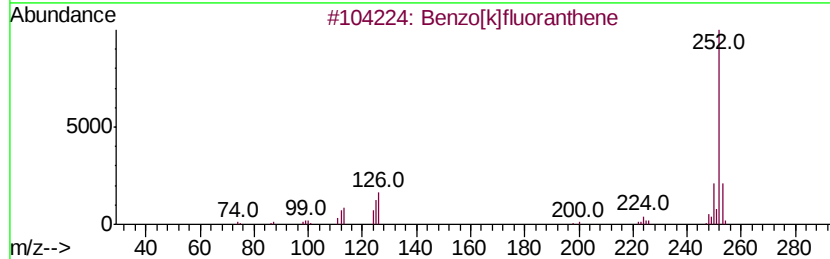
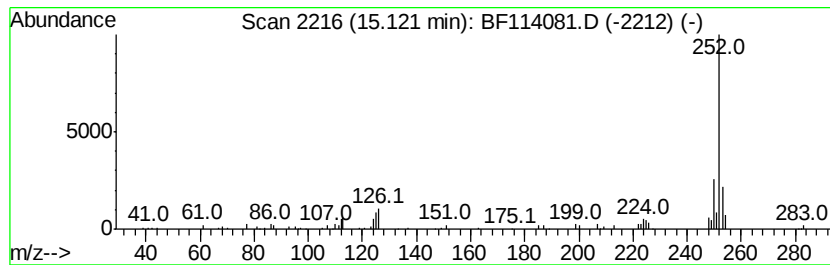
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzo[e]pyrene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.12	2.19 ng	65609	Perylene-d12	15.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
2		Benzo[e]pyrene	252	C20H12	000192-97-2	93
3		Benzo[a]acephenanthrylene	252	C20H12	000205-99-2	83
4		Benzo[a]pyrene	252	C20H12	000050-32-8	83
5		Benzo[j]fluoranthene	252	C20H12	000205-82-3	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF050219\
Data File : BF114081.D
Acq On : 2 May 2019 13:02
Operator : JU/SJ
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Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SSSI-0419-S-DH-7

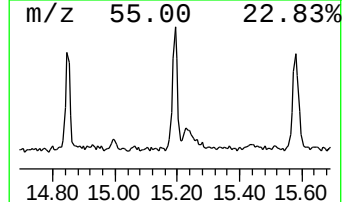
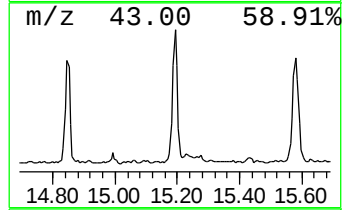
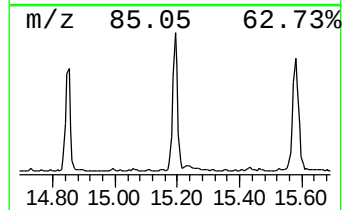
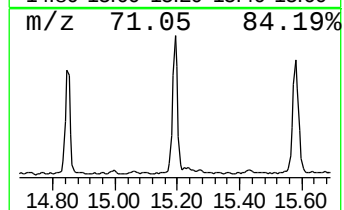
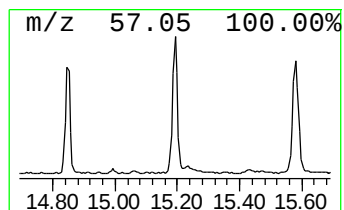
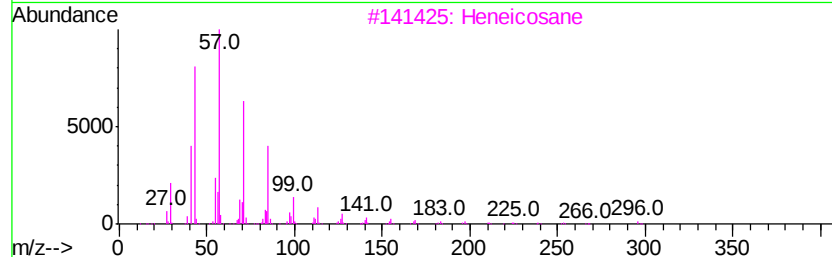
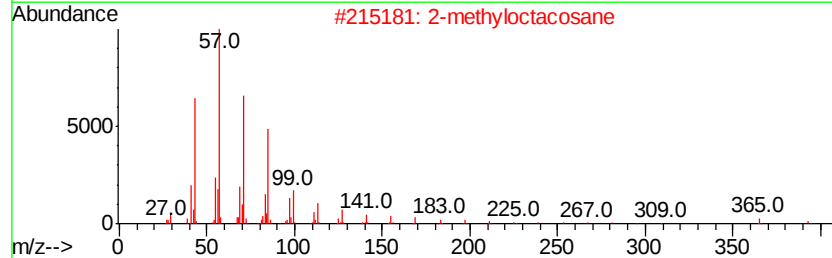
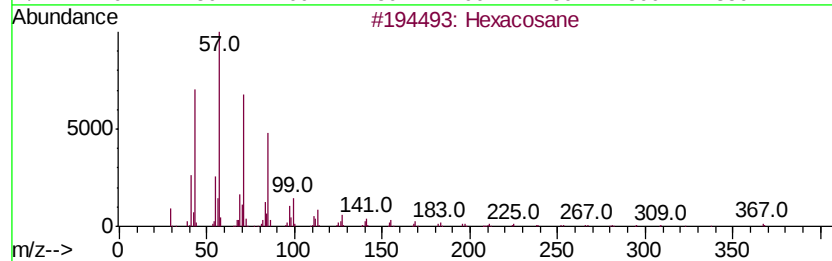
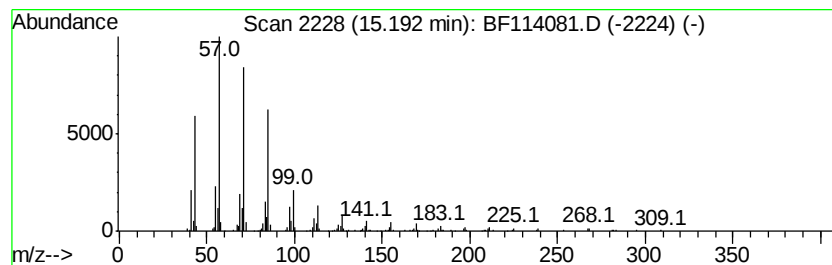
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Hexacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.19	12.34 ng	369864	Perylene-d12	15.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	91
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		triacontane	422	C30H62	000638-68-6	91
5		Heptadecane	240	C17H36	000629-78-7	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF050219\
Data File : BF114081.D
Acq On : 2 May 2019 13:02
Operator : JU/SJ
Sample : K2584-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SSSI-0419-S-DH-7

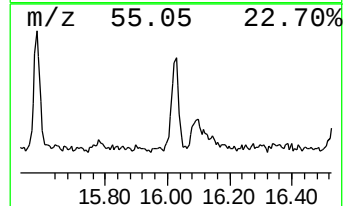
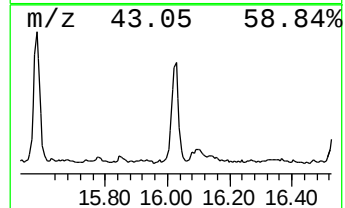
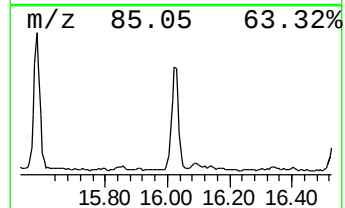
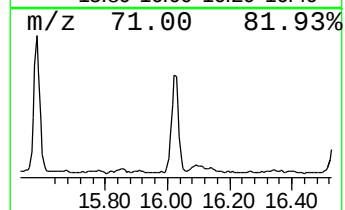
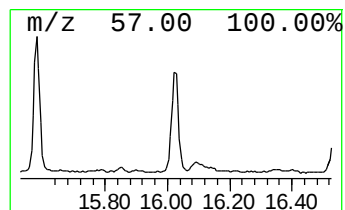
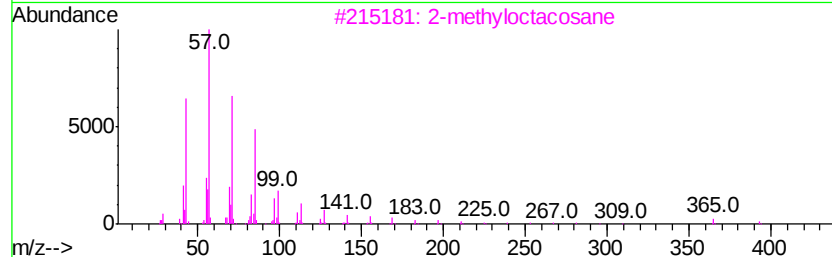
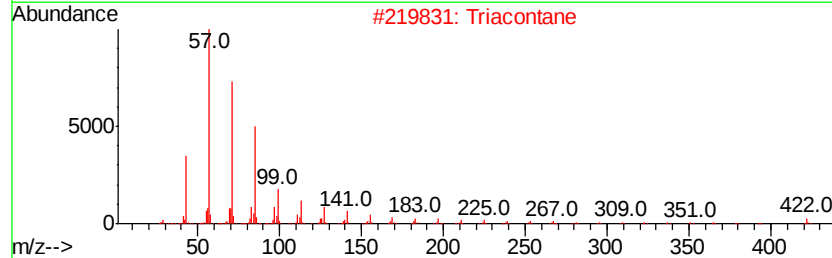
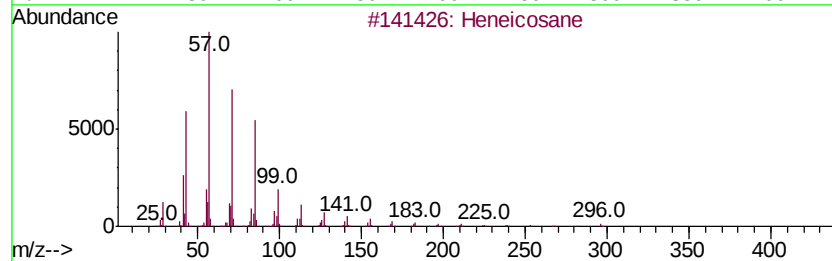
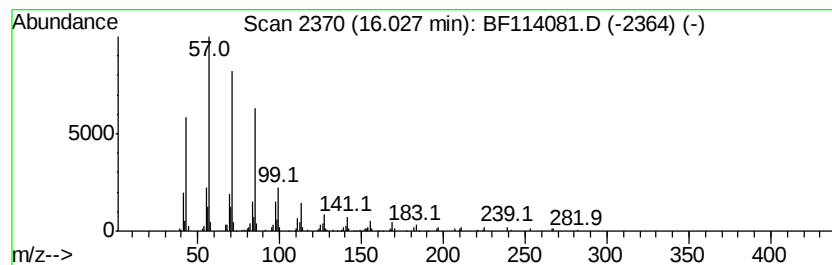
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heneicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.03	10.58 ng	317297	Perylene-d12	15.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Triacontane	422	C30H62	000638-68-6	91
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Hexacosane	366	C26H54	000630-01-3	91
5		Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF050219\
Data File : BF114081.D
Acq On : 2 May 2019 13:02
Operator : JU/SJ
Sample : K2584-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

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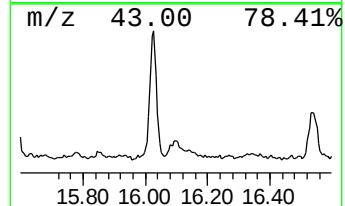
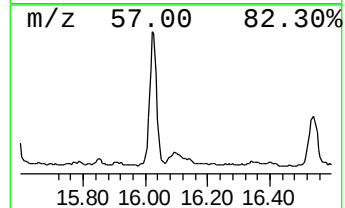
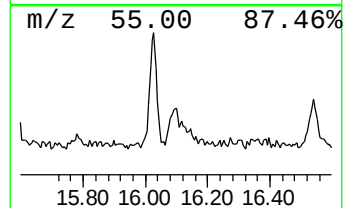
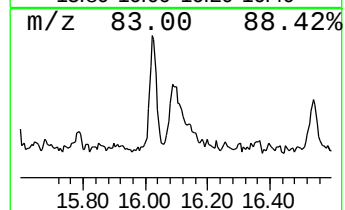
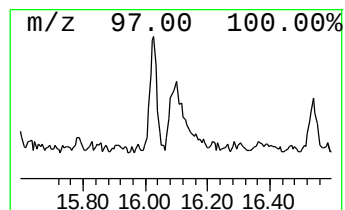
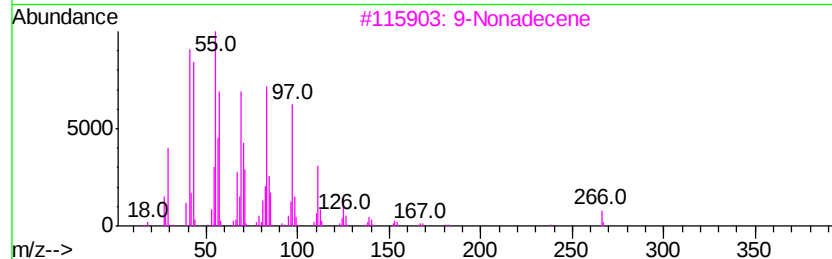
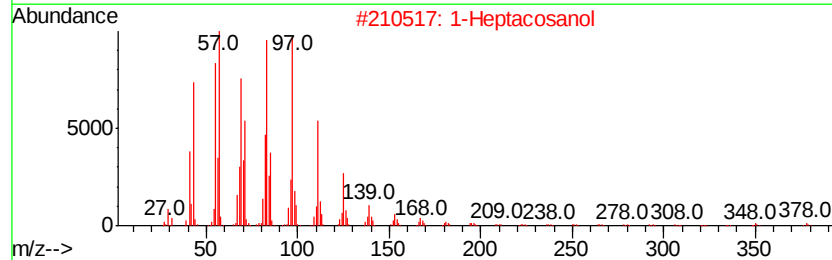
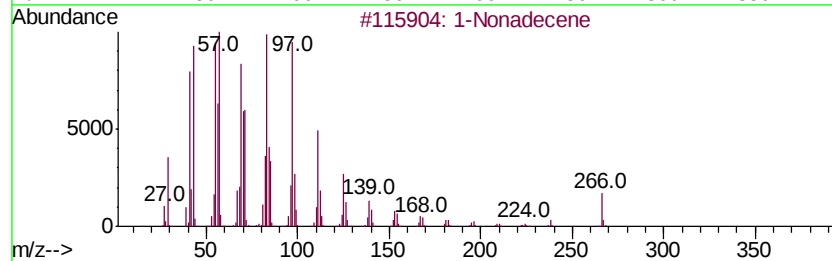
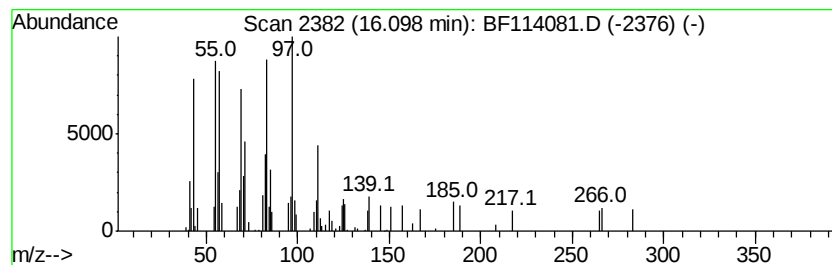
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 1-Nonadecene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.10	4.12 ng	123543	Perylene-d12	15.56

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Nonadecene	266	C19H38	018435-45-5	91
2			1-Heptacosanol	396	C27H56O	002004-39-9	86
3			9-Nonadecene	266	C19H38	031035-07-1	58
4			17-Pentatriacontene	491	C35H70	006971-40-0	49
5			Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF050219\
Data File : BF114081.D
Acq On : 2 May 2019 13:02
Operator : JU/SJ
Sample : K2584-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SSSI-0419-S-DH-7

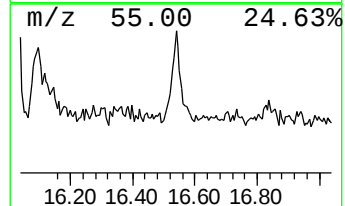
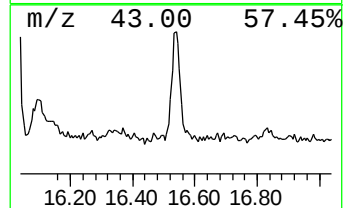
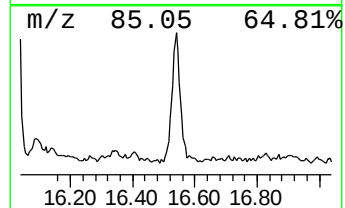
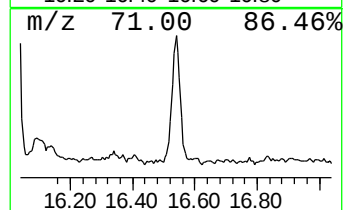
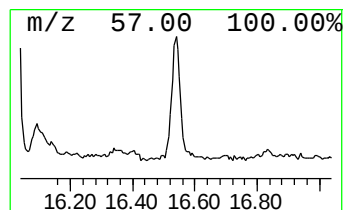
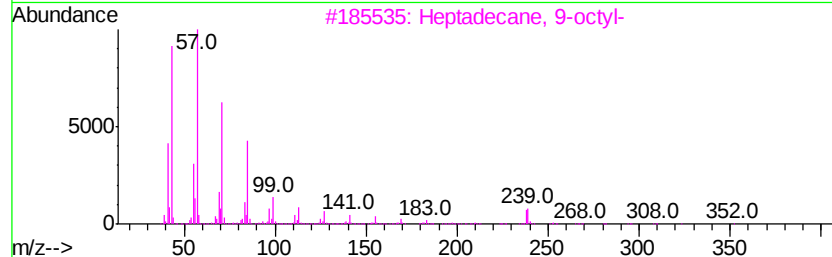
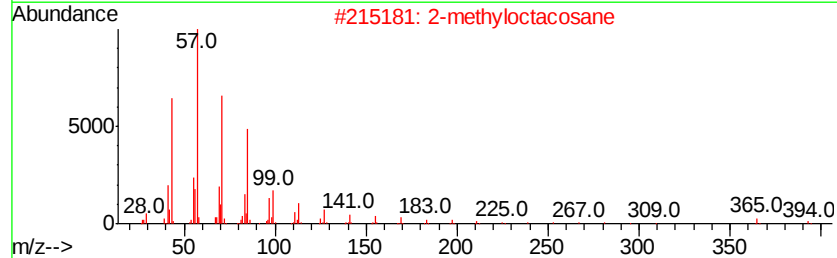
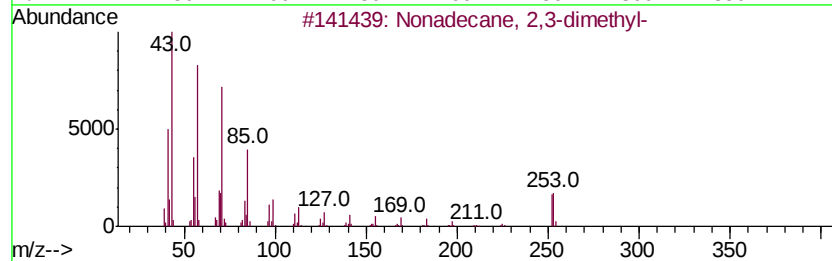
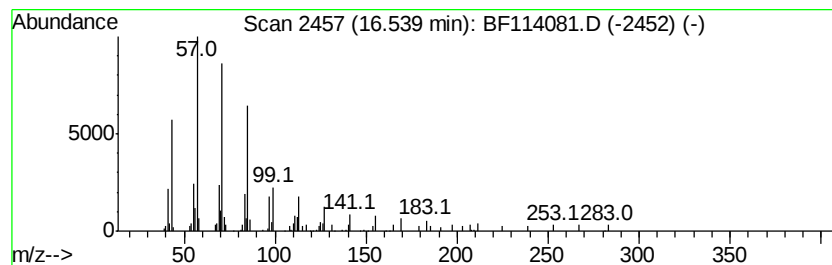
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Nonadecane, 2,3-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.54	4.67 ng	139998	Perylene-d12	15.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane, 2,3-dimethyl-	296	C21H44	075163-99-4	91
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		Hentriacontane	437	C31H64	000630-04-6	90
5		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF050219\
Data File : BF114081.D
Acq On : 2 May 2019 13:02
Operator : JU/SJ
Sample : K2584-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SSSI-0419-S-DH-7

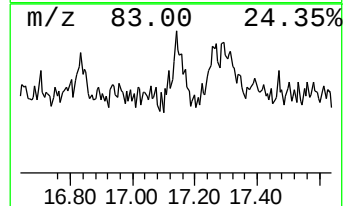
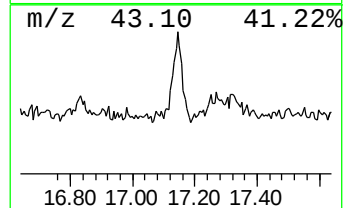
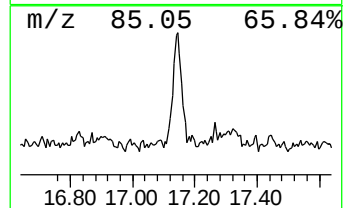
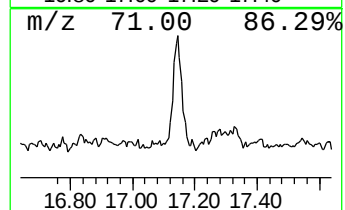
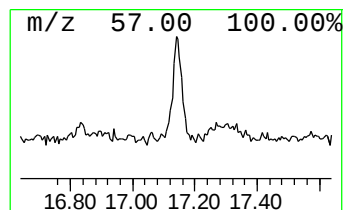
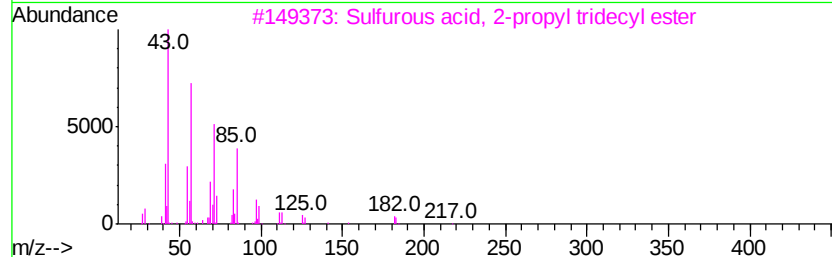
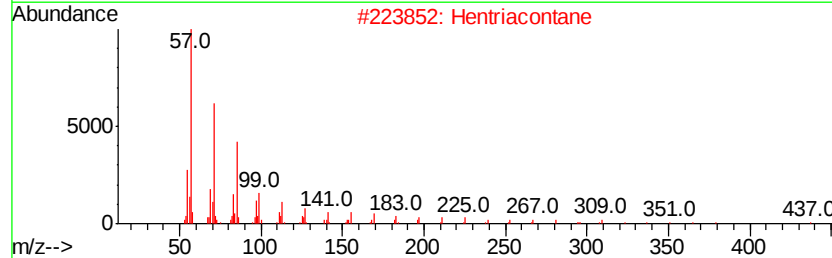
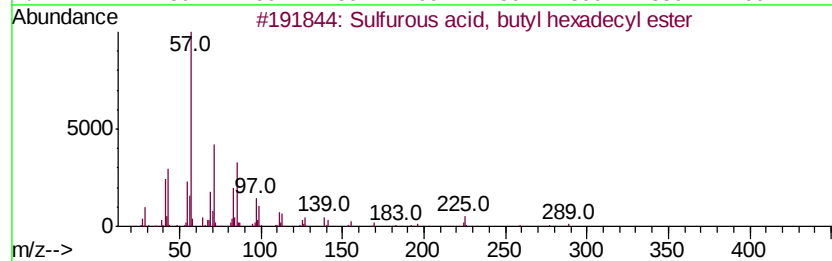
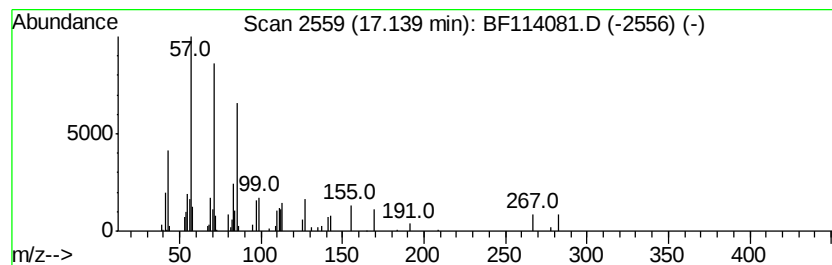
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Sulfurous acid, butyl hexad... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.14	2.28 ng	68291	Perylene-d12	15.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Sulfurous acid, butyl hexadecyl ...	362	C20H42O3S	1000309-18-3	83
2		Hentriacontane	437	C31H64	000630-04-6	78
3		Sulfurous acid, 2-propyl tridecyl...	306	C16H34O3S	1000309-12-4	72
4		2-methyloctacosane	408	C29H60	1000376-72-8	72
5		Tetradecane	198	C14H30	000629-59-4	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF050219\
Data File : BF114081.D
Acq On : 2 May 2019 13:02
Operator : JU/SJ
Sample : K2584-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SSSI-0419-S-DH-7

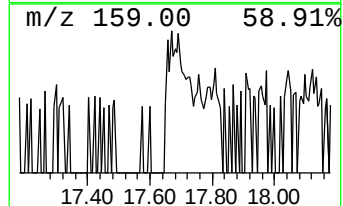
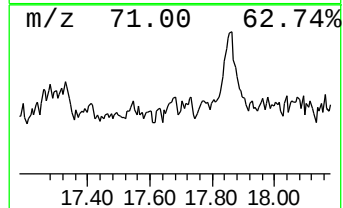
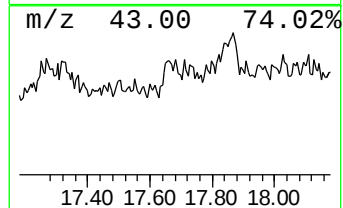
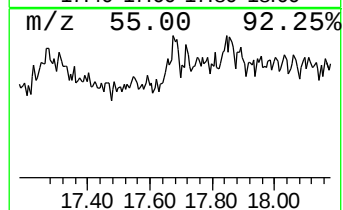
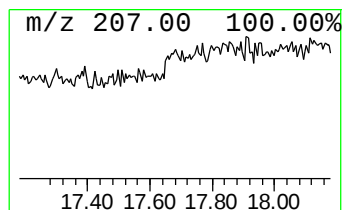
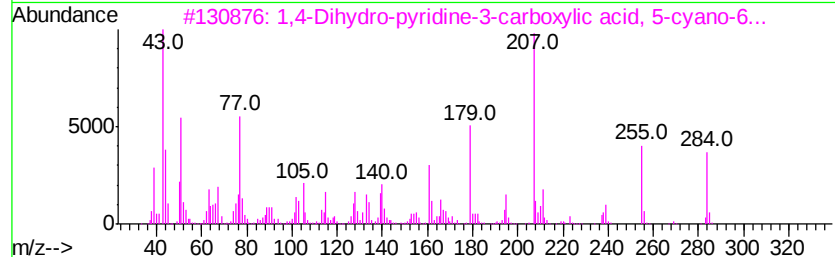
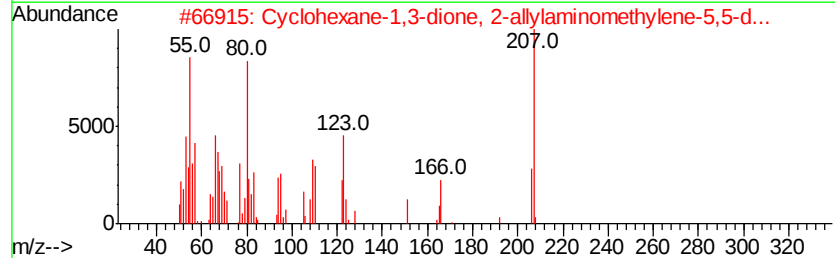
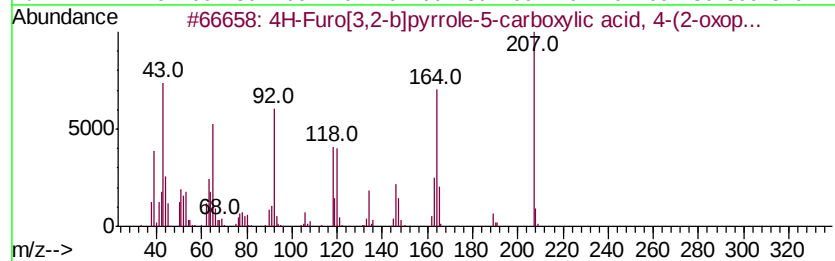
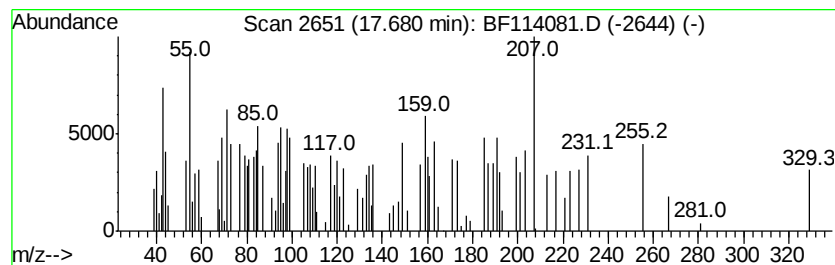
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF042219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown17.68 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.68	3.06 ng	91643	Perylene-d12	15.56

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Furo[3,2-b]pyrrole-5-carboxyl...	207	C10H9NO4	1000318-88-7	10
2			Cyclohexane-1,3-dione, 2-allylam...	207	C12H17NO2	104926-37-6	10
3			1,4-Dihydro-pyridine-3-carboxyli...	284	C16H16N2O3	1000259-17-8	10
4			1,2,5-Oxadiazol-3-amine, 4-(4-me...	207	C9H9N3O3	1000351-72-7	10
5			Indole-2-one, 2,3-dihydro-N-hydr...	207	C11H13NO3	1000129-52-1	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF050219\
 Data File : BF114081.D
 Acq On : 2 May 2019 13:02
 Operator : JU/SJ
 Sample : K2584-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSSI-0419-S-DH-7

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
1,3,5-Cycloheptat...	3.98	5.7	ng	123312	1	6.89	436888	20.0
unknown6.66	6.66	152.0	ng	3321070	1	6.89	436888	20.0
n-Hexadecanoic acid	11.93	12.7	ng	389740	4	11.41	614080	20.0
Octadecanoic acid	12.72	20.8	ng	639892	4	11.41	614080	20.0
Behenic alcohol	13.89	11.8	ng	416400	5	14.06	707848	20.0
Octacosane	14.22	2.5	ng	90234	5	14.06	707848	20.0
Triacontane	14.53	7.9	ng	278433	5	14.06	707848	20.0
Eicosane	14.85	9.1	ng	273933	6	15.56	599591	20.0
Benzo[e]pyrene	15.12	2.2	ng	65609	6	15.56	599591	20.0
Hexacosane	15.19	12.3	ng	369864	6	15.56	599591	20.0
Heneicosane	16.03	10.6	ng	317297	6	15.56	599591	20.0
1-Nonadecene	16.10	4.1	ng	123543	6	15.56	599591	20.0
Nonadecane, 2,3-d...	16.54	4.7	ng	139998	6	15.56	599591	20.0
Sulfurous acid, b...	17.14	2.3	ng	68291	6	15.56	599591	20.0
unknown17.68	17.68	3.1	ng	91643	6	15.56	599591	20.0