

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012615\
 Data File : BF077082.D
 Acq On : 27 Jan 2015 9:59
 Operator : TP/IZ
 Sample : G1160-03
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC012215-SOIL

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:\HPCHEM1\BNA_F\METHODS\8270-BF011415.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	3.830	237	240	247	rBV	37930	85293	1.04%	0.364%
2	4.801	322	325	337	rBV	250221	496438	6.07%	2.119%
3	7.179	530	533	538	rBV	262180	499720	6.11%	2.133%
4	7.259	538	540	547	rVB	334491	546415	6.68%	2.332%
5	7.762	581	584	589	rBV	82247	134929	1.65%	0.576%
6	8.082	609	612	617	rBV	227093	384254	4.70%	1.640%
7	9.065	695	698	706	rBV	205404	319679	3.91%	1.364%
8	10.676	835	839	842	rBV	126726	192012	2.35%	0.819%
9	13.328	1068	1071	1075	rBV	365802	576536	7.05%	2.460%
10	14.802	1196	1200	1204	rVB2	126561	210315	2.57%	0.898%
11	16.471	1343	1346	1348	rBV	82106	118454	1.45%	0.506%
12	16.517	1348	1350	1353	rVV	272104	380907	4.66%	1.626%
13	17.637	1445	1448	1450	rBV	171012	233388	2.85%	0.996%
14	18.460	1516	1520	1522	rVB	1510823	2057768	25.15%	8.782%
15	18.643	1532	1536	1540	rBV2	68868	125967	1.54%	0.538%
16	19.363	1592	1599	1600	rBV	4692721	8181056	100.00%	34.913%
17	19.386	1600	1601	1604	rVB	3813377	5220896	63.82%	22.281%
18	19.500	1609	1611	1614	rVV	902745	976629	11.94%	4.168%
19	19.877	1641	1644	1647	rBV	555784	637717	7.80%	2.721%
20	20.312	1679	1682	1684	rBV2	64834	93898	1.15%	0.401%
21	20.426	1690	1692	1696	rBV2	186690	287649	3.52%	1.228%
22	20.494	1696	1698	1701	rVV	78487	118909	1.45%	0.507%
23	20.677	1710	1714	1717	rVV4	84000	154021	1.88%	0.657%
24	20.952	1735	1738	1741	rBV3	71012	137376	1.68%	0.586%
25	21.135	1751	1754	1756	rBV2	226403	423902	5.18%	1.809%
26	21.283	1763	1767	1770	rBV2	160043	304166	3.72%	1.298%
27	22.037	1831	1833	1839	rVB2	65271	137369	1.68%	0.586%
28	22.780	1896	1898	1901	rVB	184412	218586	2.67%	0.933%
29	23.375	1947	1950	1955	rVB5	74106	178315	2.18%	0.761%

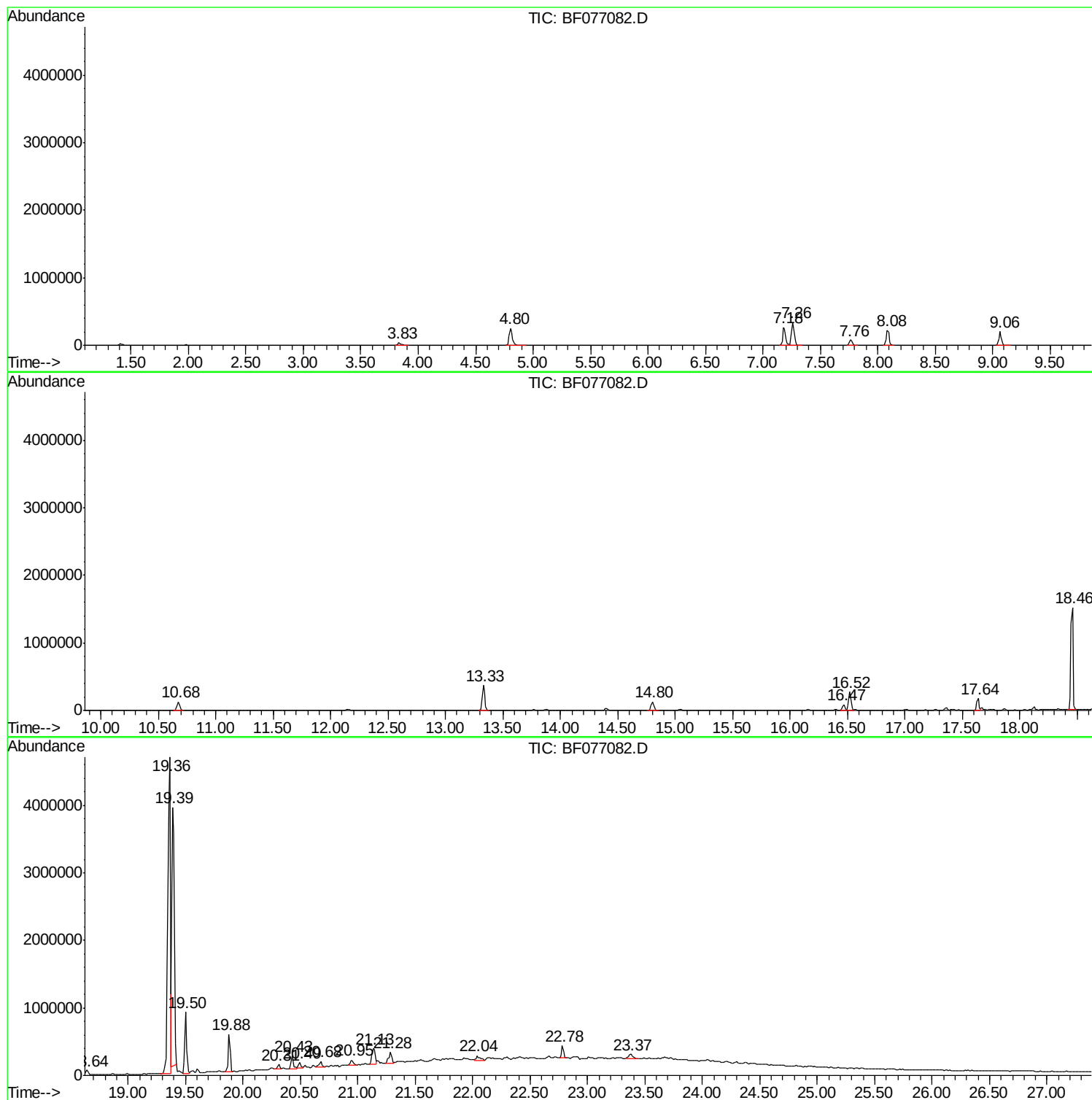
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Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF011415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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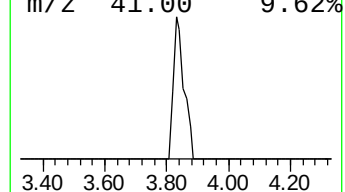
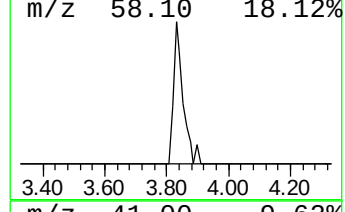
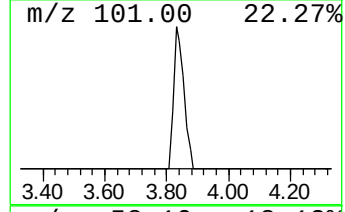
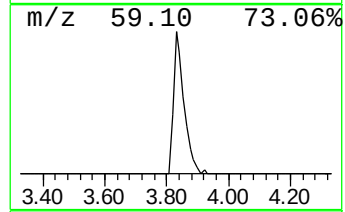
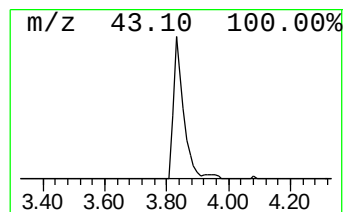
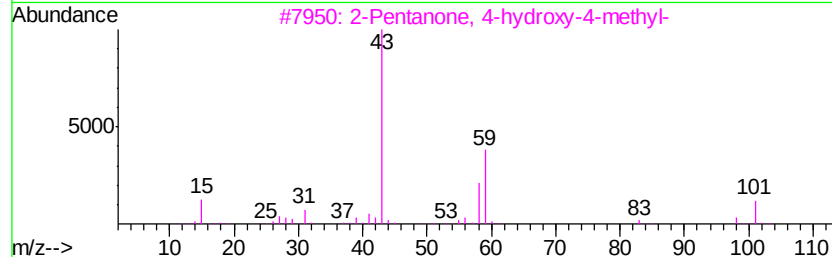
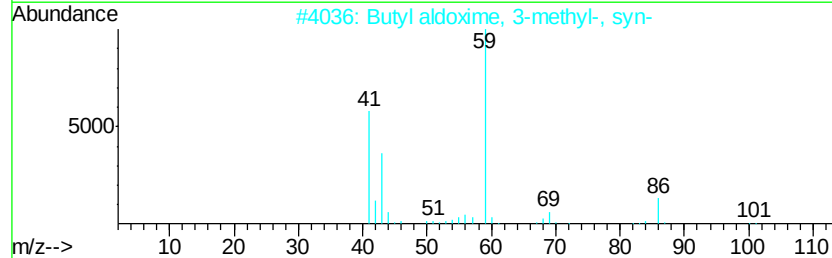
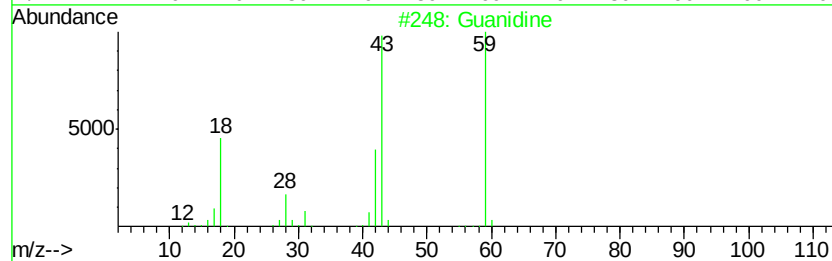
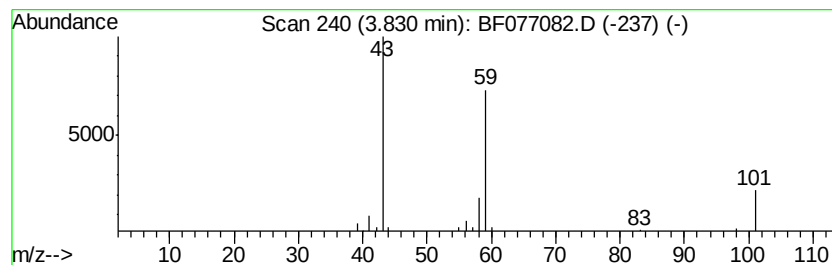
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF011415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown3.83 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.83	12.64 ng	85293	1,4-Dichlorobenzene-d4	7.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Guanidine	59	CH5N3	000113-00-8	43
2		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	37
3		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	33
4		3-Hexanol	102	C6H14O	000623-37-0	25
5		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	23



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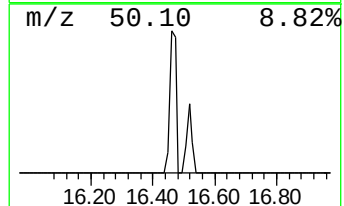
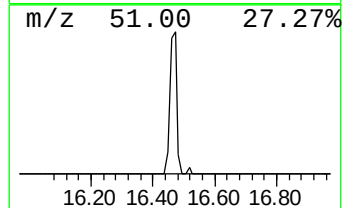
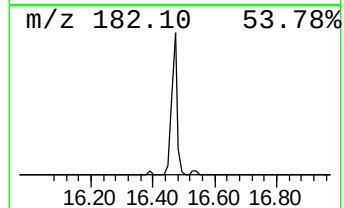
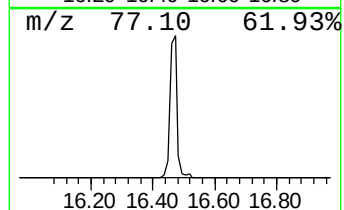
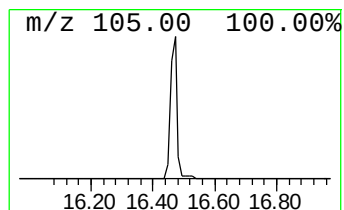
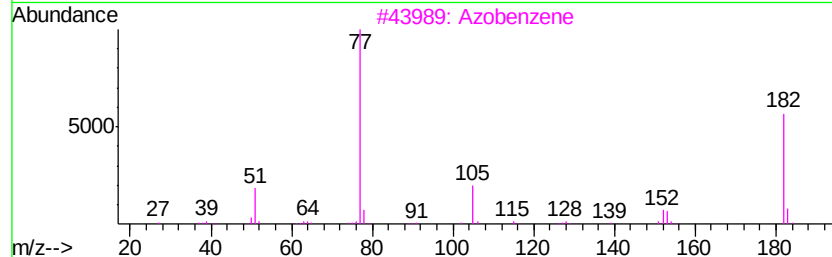
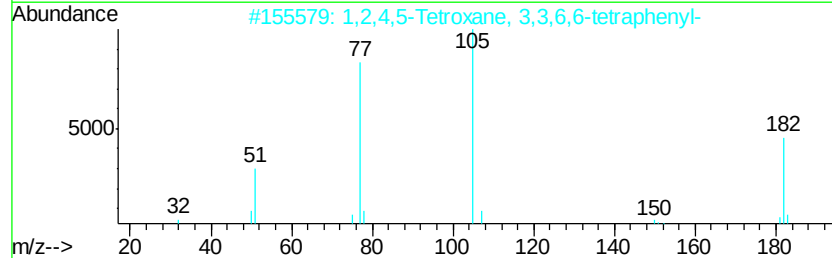
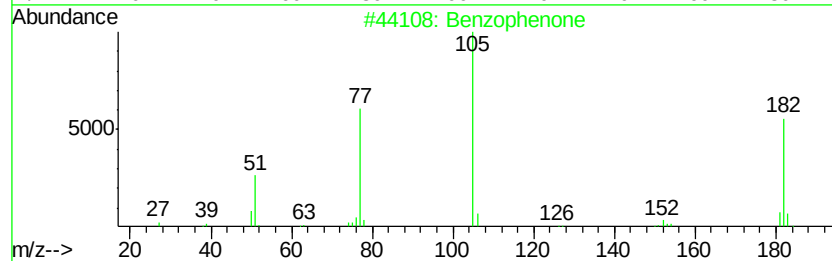
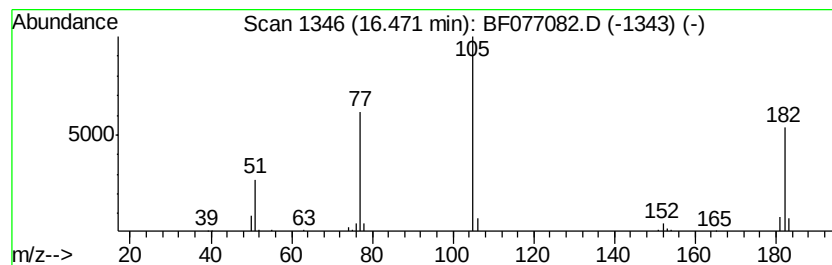
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 Benzophenone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.47	10.15 ng	118454	Phenanthrene-d10	17.64

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzophenone	182	C13H10O	000119-61-9	96
2		1,2,4,5-Tetroxane, 3,3,6,6-tetra...	396	C26H20O4	016204-36-7	78
3		Azobenzene	182	C12H10N2	000103-33-3	59
4		Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	47
5		Benzoylformic acid	150	C8H6O3	000611-73-4	47



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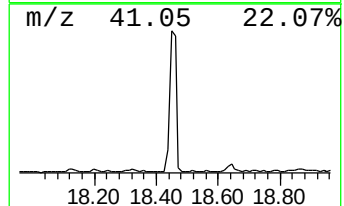
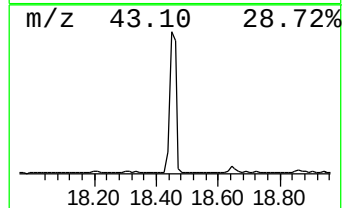
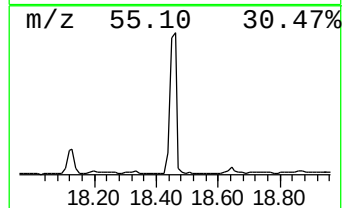
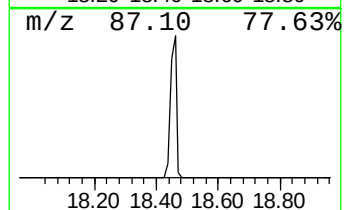
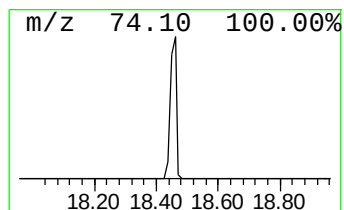
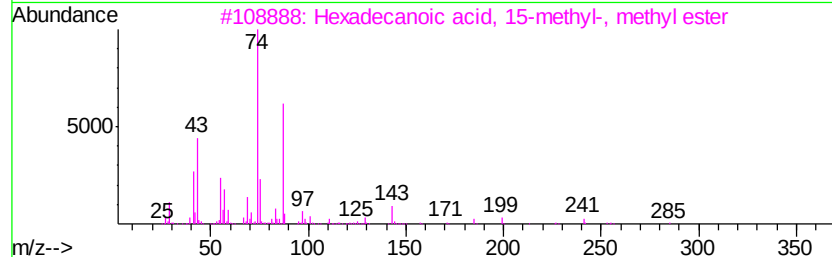
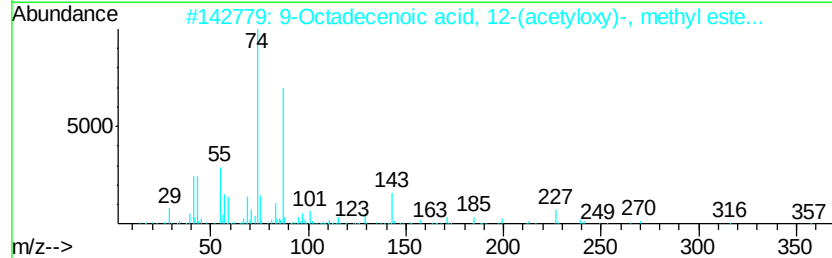
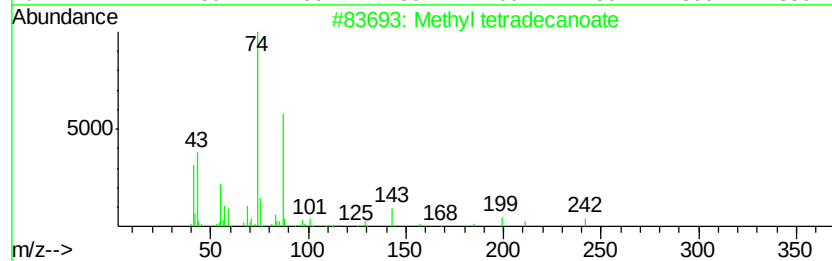
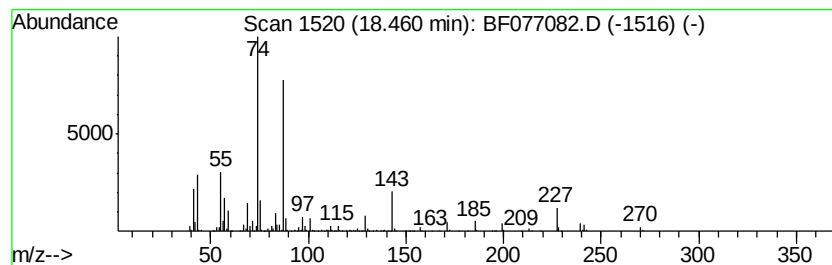
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 4 Methyl tetradecanoate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.46	176.34 ng	2057770	Phenanthrene-d10	17.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methyl tetradecanoate	242	C15H30O2	000124-10-7	91
2		9-Octadecenoic acid, 12-(acetylo...	354	C21H38O4	000140-03-4	87
3		Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	83
4		Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	83
5		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	81



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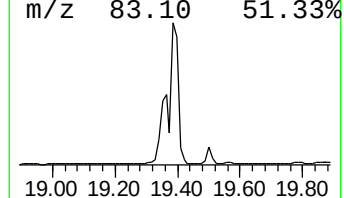
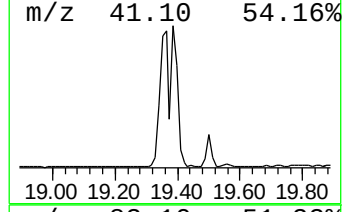
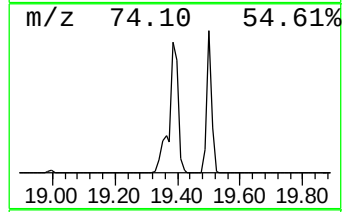
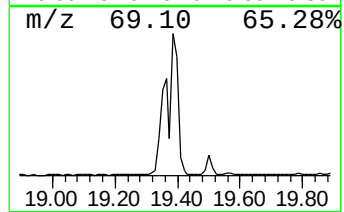
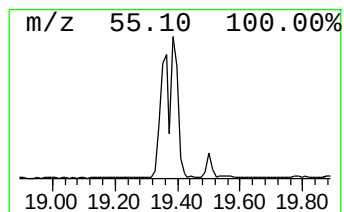
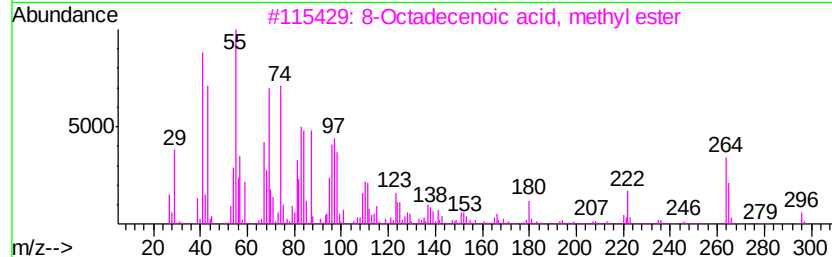
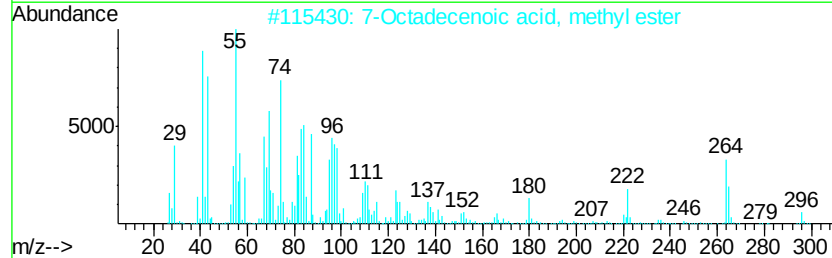
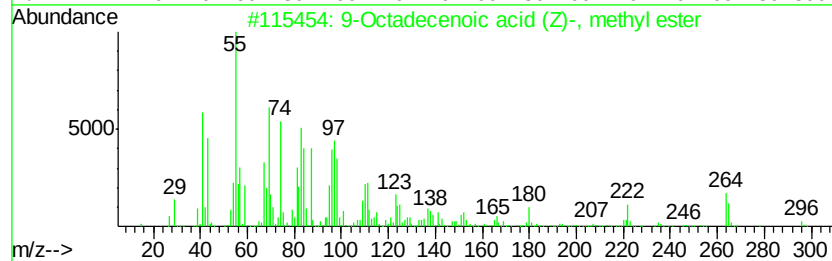
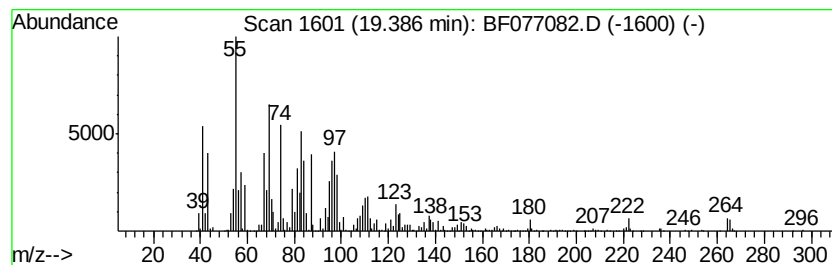
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 9-Octadecenoic acid (Z)-, m... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.39	447.40 ng	5220900	Phenanthrene-d10	17.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	99
2		7-Octadecenoic acid, methyl ester	296	C19H36O2	057396-98-2	98
3		8-Octadecenoic acid, methyl ester	296	C19H36O2	002345-29-1	97
4		9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	94
5		12-Octadecenoic acid, methyl ester	296	C19H36O2	056554-46-2	91



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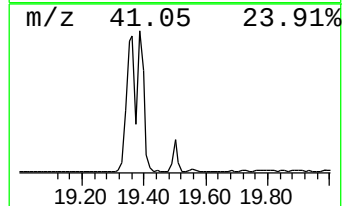
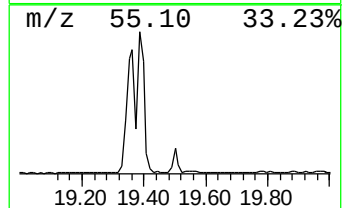
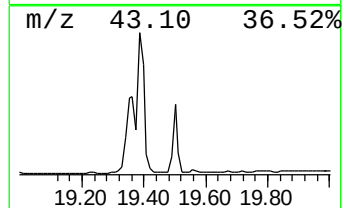
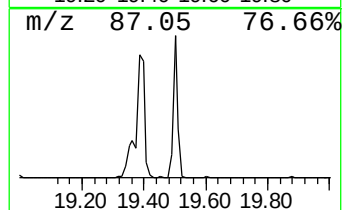
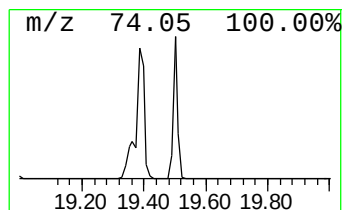
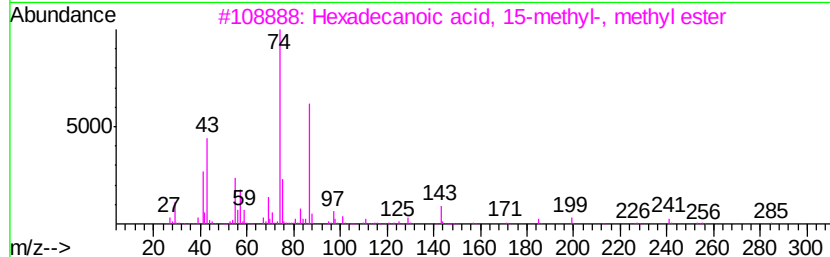
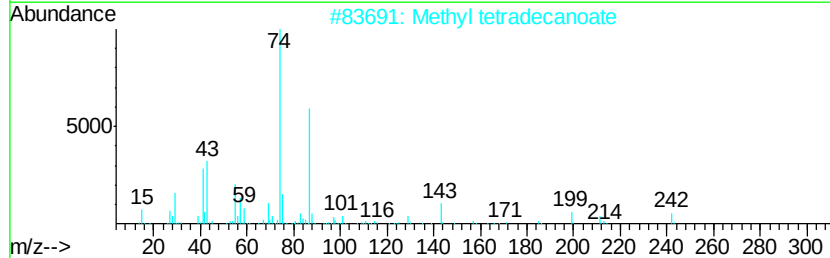
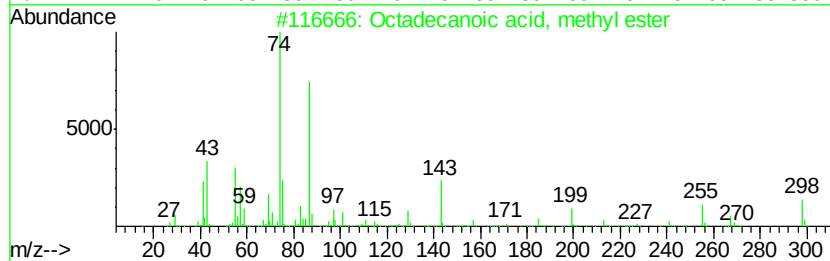
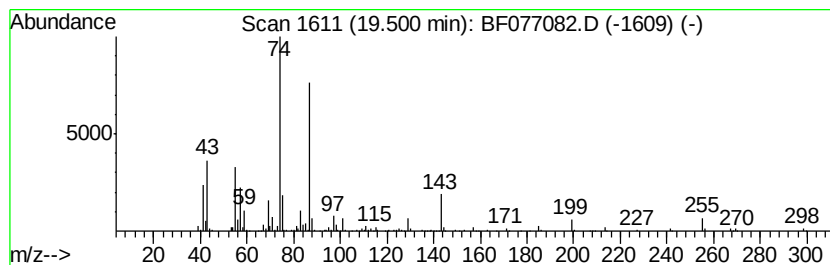
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 6 Octadecanoic acid, methyl e... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.50	46.08 ng	976629	Chrysene-d12	21.15	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid, methyl ester	298	C19H38O2	000112-61-8	98
2	Methyl tetradecanoate	242	C15H30O2	000124-10-7	91
3	Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	91
4	Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	91
5	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	90



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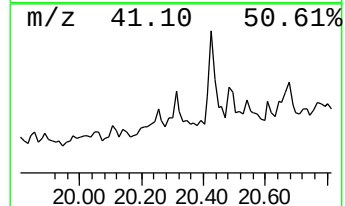
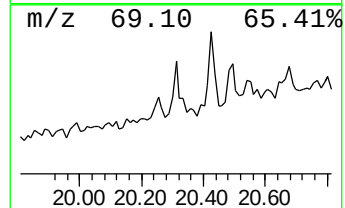
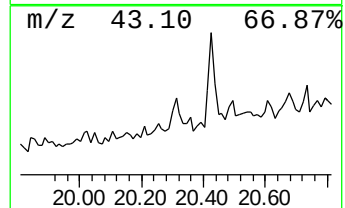
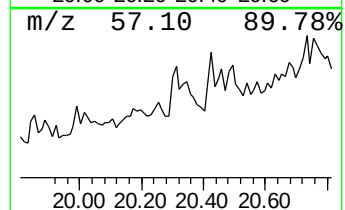
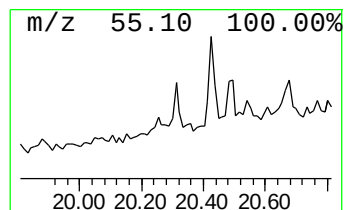
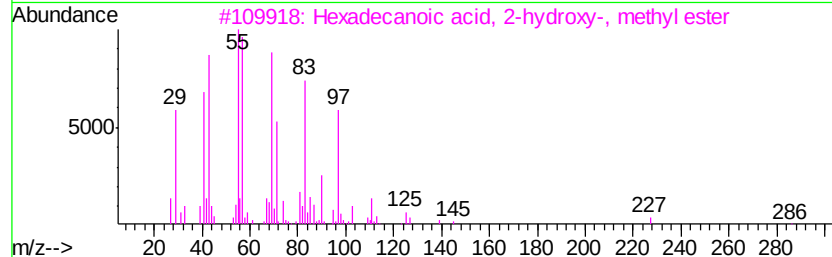
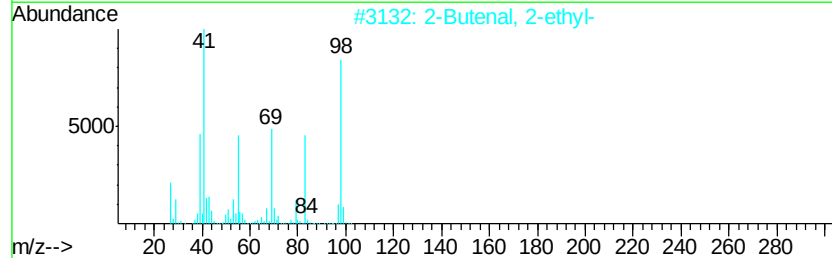
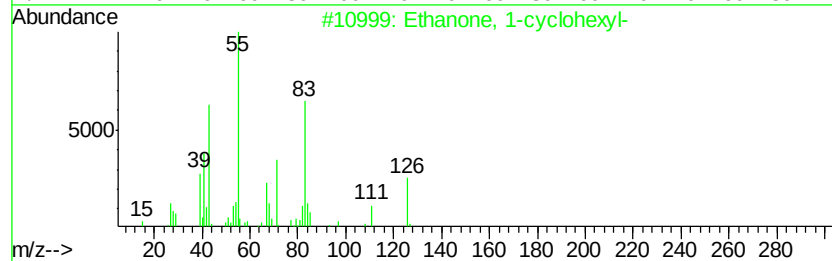
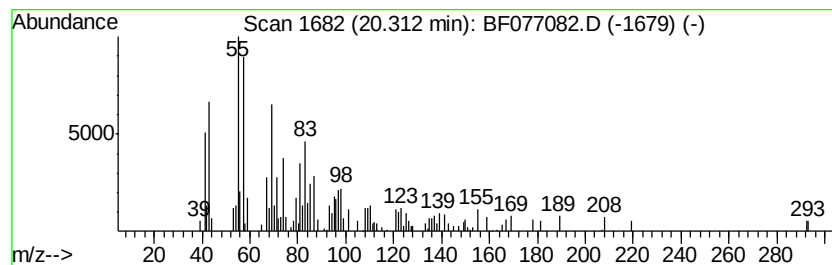
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Ethanone, 1-cyclohexyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.		
20.31	4.43 ng	93898	Chrysene-d12	21.15		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanone, 1-cyclohexyl-	126	C8H14O	000823-76-7	60	
2	2-Butenal, 2-ethyl-	98	C6H10O	019780-25-7	38	
3	Hexadecanoic acid, 2-hydroxy-, m...	286	C17H34O3	016742-51-1	35	
4	3-Hexen-2-one, 3-cyclohexyl-4-et...	208	C14H24O	1000150-19-1	30	
5	Oxiraneundecanoic acid, 3-pentyl...	312	C19H36O3	038520-30-8	30	



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012615\
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Acq On : 27 Jan 2015 9:59
Operator : TP/IZ
Sample : G1160-03
Misc :
ALS Vial : 36 Sample Multiplier: 1

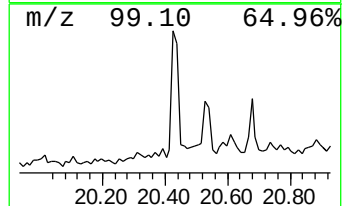
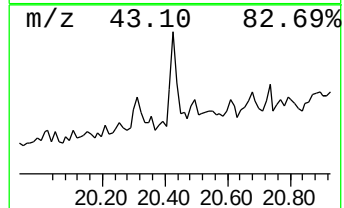
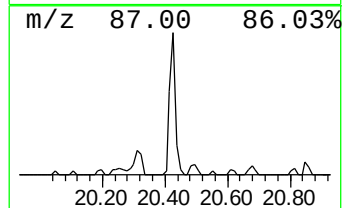
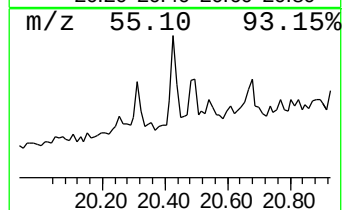
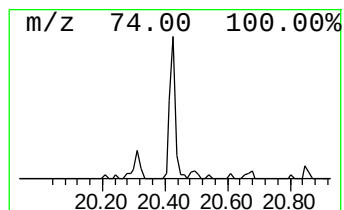
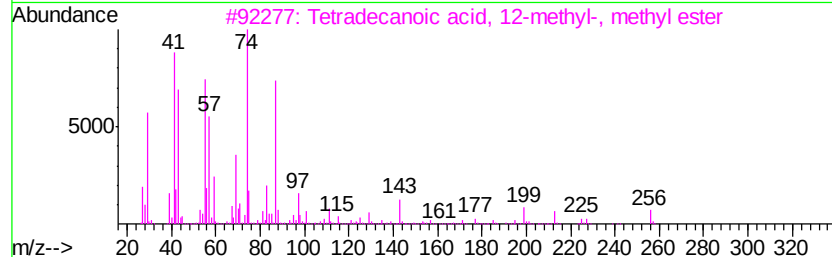
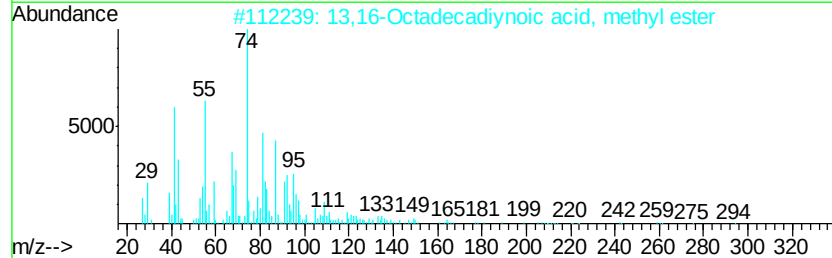
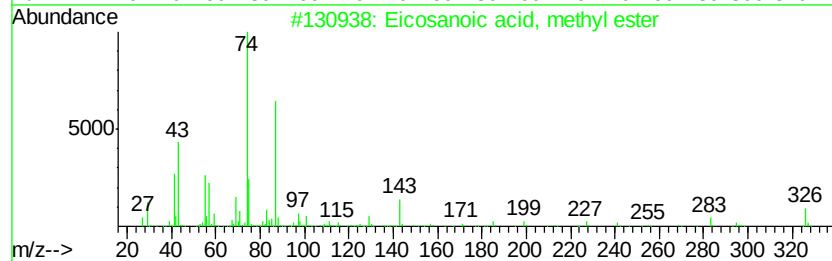
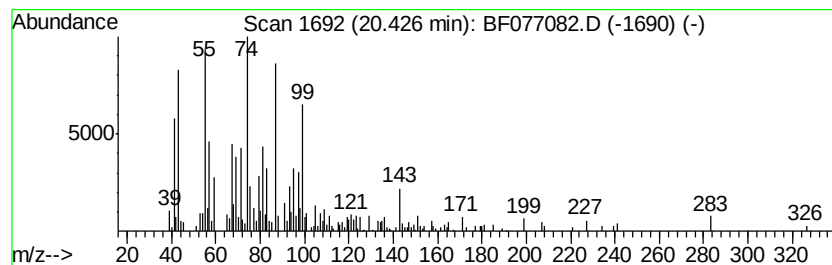
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Eicosanoic acid, methyl ester Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.		
20.43	13.57 ng	287649	Chrysene-d12	21.15		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosanoic acid, methyl ester	326	C21H42O2	001120-28-1	64	
2	13,16-Octadecadiynoic acid, meth...	290	C19H30O2	056846-98-1	49	
3	Tetradecanoic acid, 12-methyl-, ...	256	C16H32O2	005129-66-8	43	
4	Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	38	
5	Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	38	



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012615\
Data File : BF077082.D
Acq On : 27 Jan 2015 9:59
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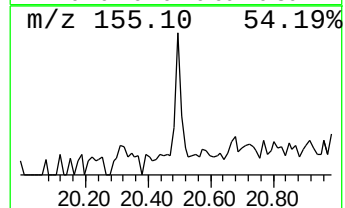
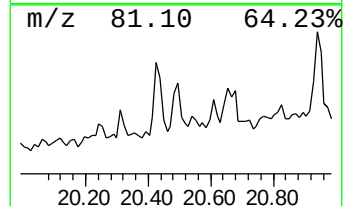
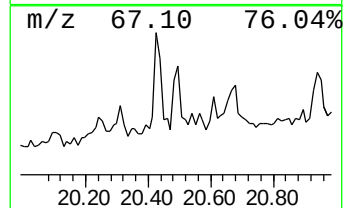
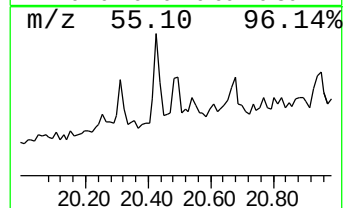
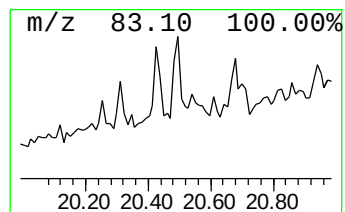
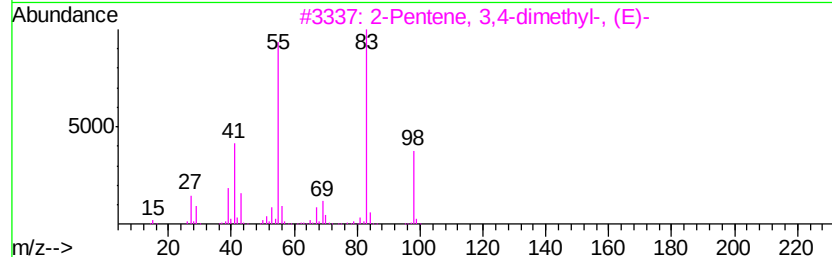
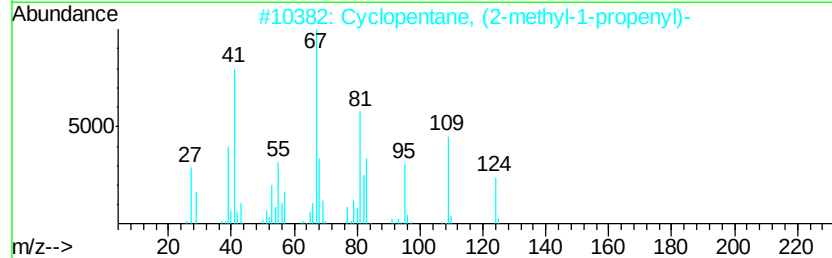
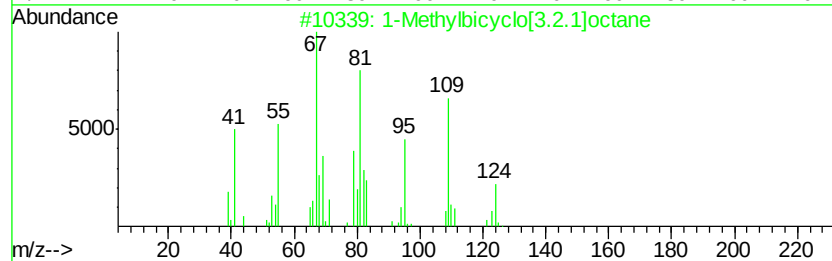
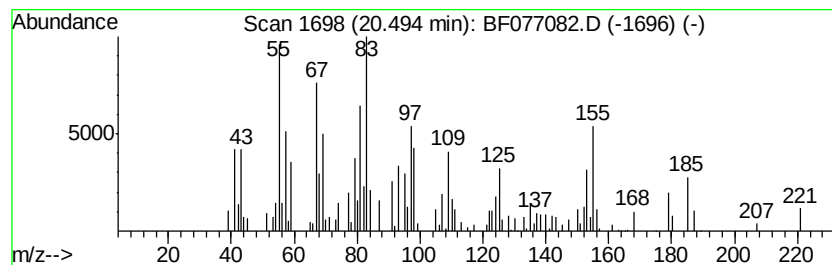
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 unknown20.49 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.49	5.61 ng	118909	Chrysene-d12	21.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Methylbicyclo[3.2.1]octane	124	C9H16	119972-41-7	35
2			Cyclopentane, (2-methyl-1-propen...	124	C9H16	053366-57-7	22
3			2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	18
4			2-Pentene, 2,3-dimethyl-	98	C7H14	010574-37-5	18
5			2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	18



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012615\
Data File : BF077082.D
Acq On : 27 Jan 2015 9:59
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ClientSampleId :
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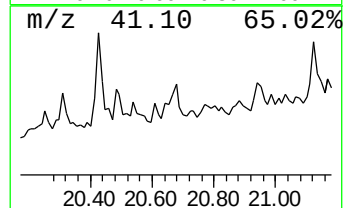
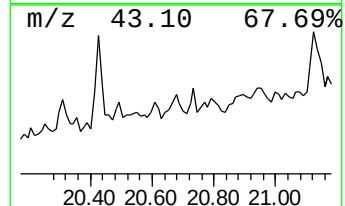
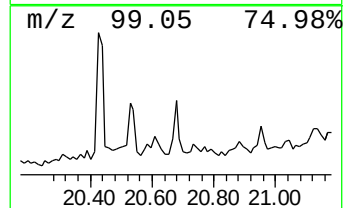
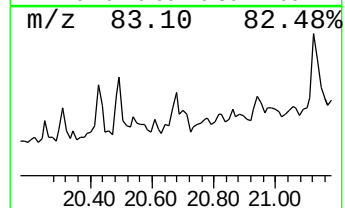
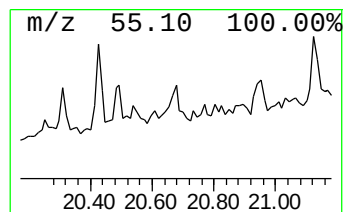
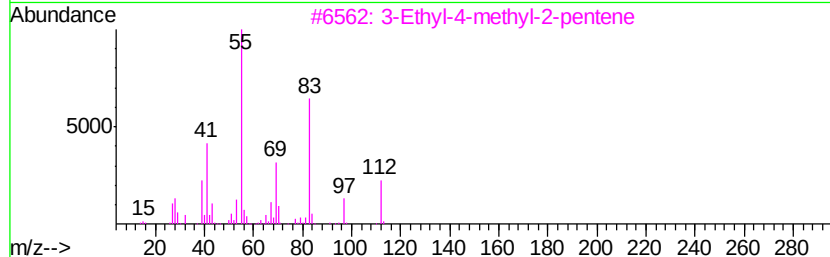
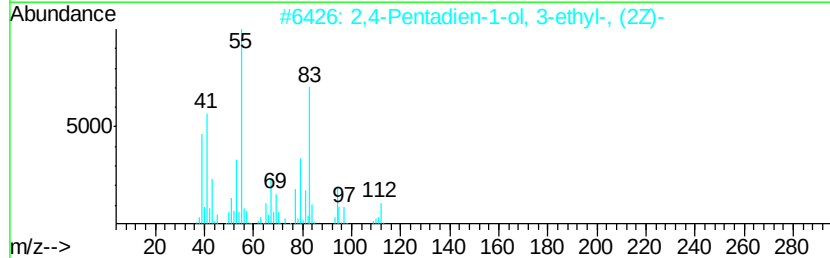
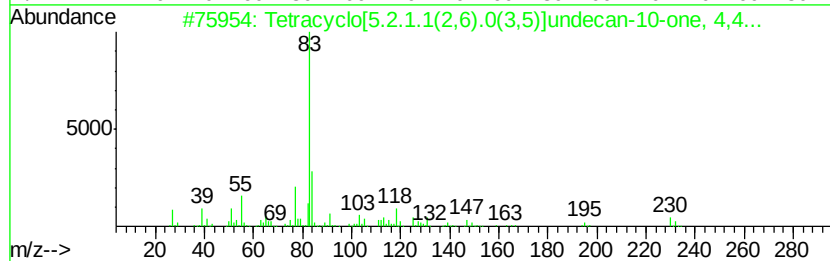
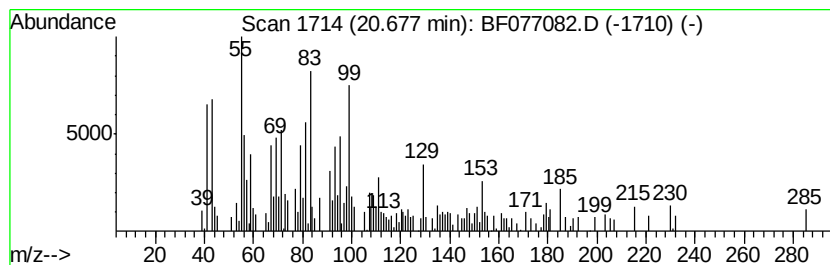
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 unknown20.68 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.68	7.27 ng	154021	Chrysene-d12	21.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracyclo[5.2.1.1(2,6).0(3,5)]u...	230	C11H12Cl2O	1000187-02-5	15
2		2,4-Pentadien-1-ol, 3-ethyl-, (2Z)-	112	C7H12O	1000142-17-6	15
3		3-Ethyl-4-methyl-2-pentene	112	C8H16	019780-68-8	15
4		Anhydro-6,8-dimethyl-1-[p-methox...	285	C15H15N3O3	1000214-73-0	11
5		3,4-Dimethyl-2-pentene(c,t)	98	C7H14	1000118-16-5	11



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012615\
Data File : BF077082.D
Acq On : 27 Jan 2015 9:59
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Misc :
ALS Vial : 36 Sample Multiplier: 1

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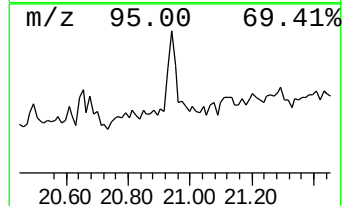
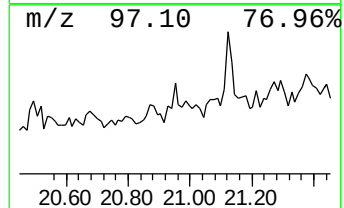
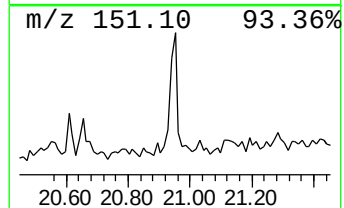
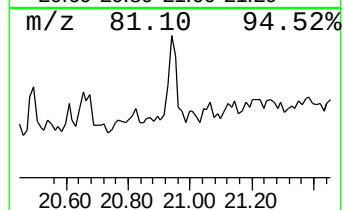
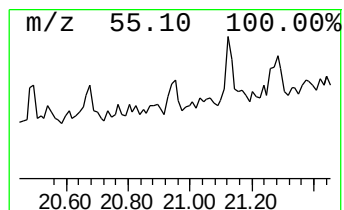
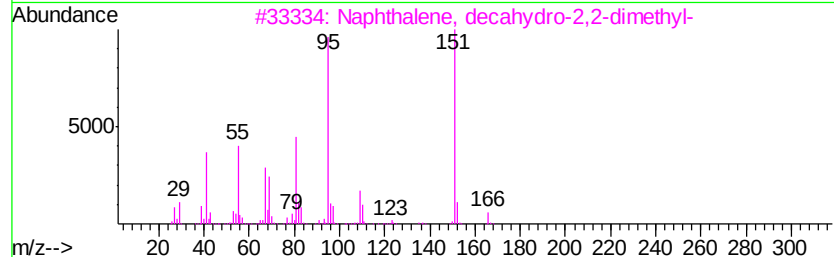
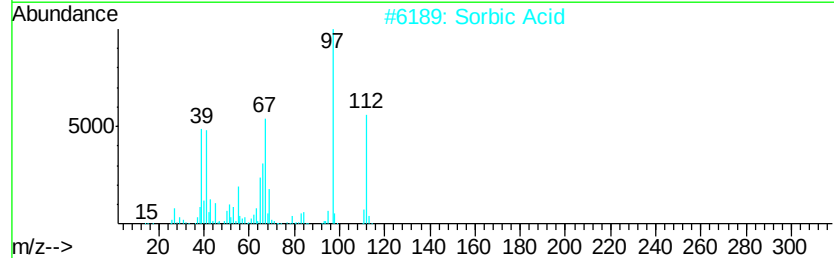
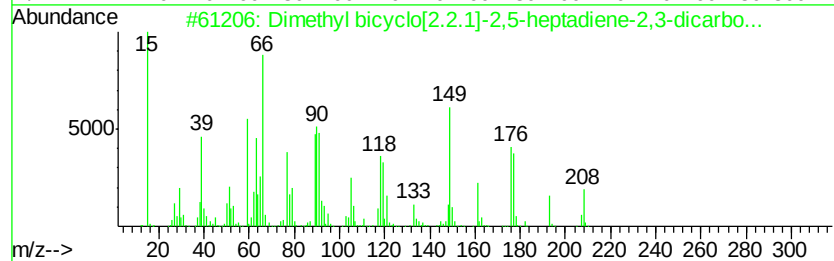
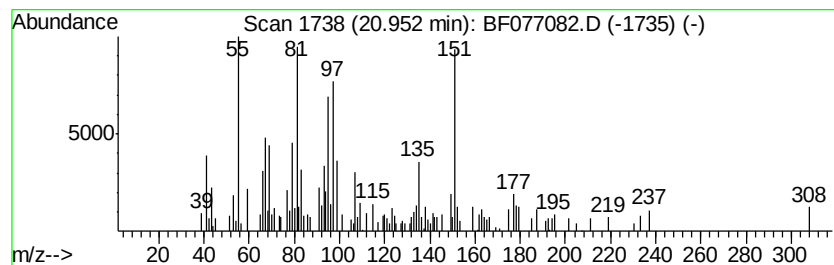
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 unknown20.95 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.95	6.48 ng	137376	Chrysene-d12	21.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dimethyl bicyclo[2.2.1]-2,5-hept...	208	C11H12O4	000947-57-9	25
2		Sorbic Acid	112	C6H8O2	000110-44-1	22
3		Naphthalene, decahydro-2,2-dimet...	166	C12H22	031124-85-3	20
4		cis-1-Decalone	152	C10H16O	1000145-07-6	11
5		Piperonylamine	151	C8H9NO2	002620-50-0	11



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012615\
Data File : BF077082.D
Acq On : 27 Jan 2015 9:59
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Misc :
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ClientSampleId :
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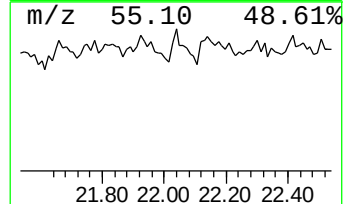
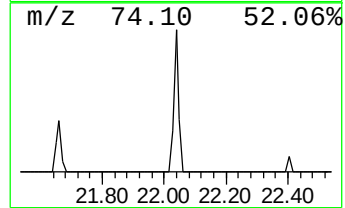
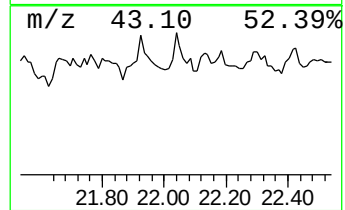
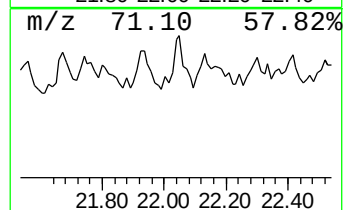
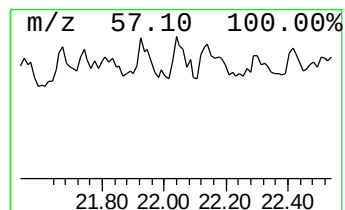
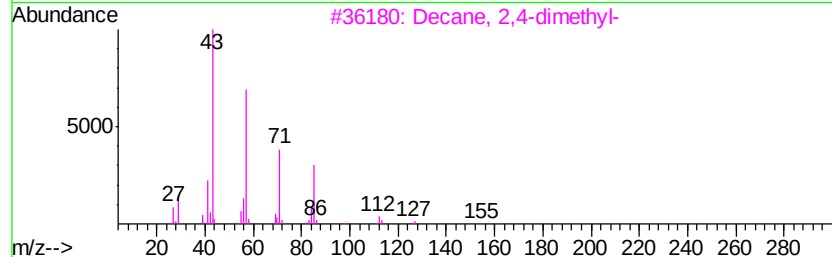
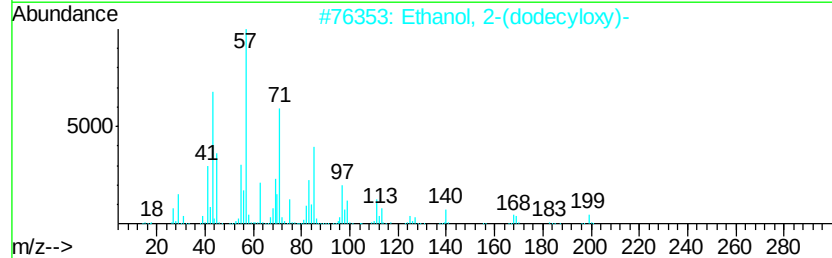
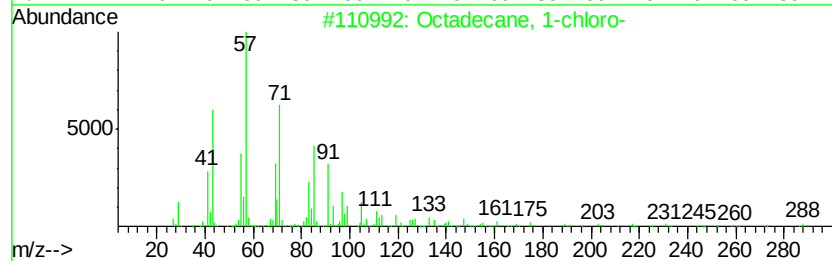
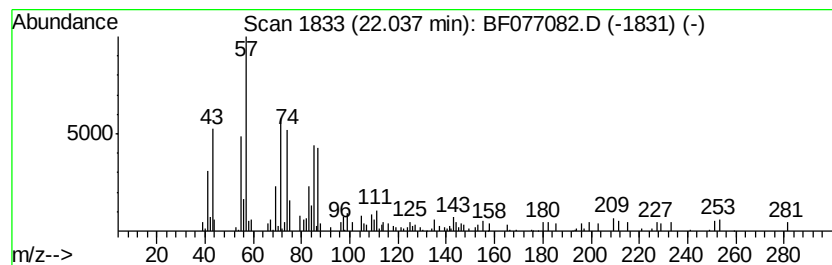
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 unknown22.04 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.04	12.57 ng	137369	Perylene-d12	22.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	35
2		Ethanol, 2-(dodecyloxy)-	230	C14H30O2	004536-30-5	35
3		Decane, 2,4-dimethyl-	170	C12H26	002801-84-5	27
4		Undecane, 4,7-dimethyl-	184	C13H28	017301-32-5	27
5		Methyl tetradecanoate	242	C15H30O2	000124-10-7	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012615\
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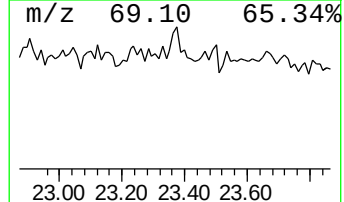
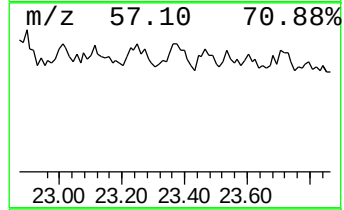
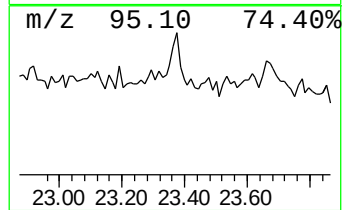
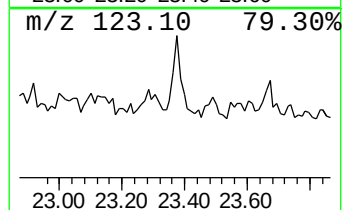
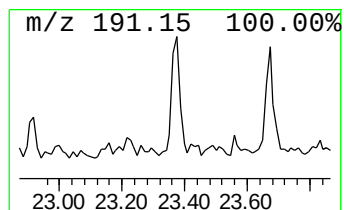
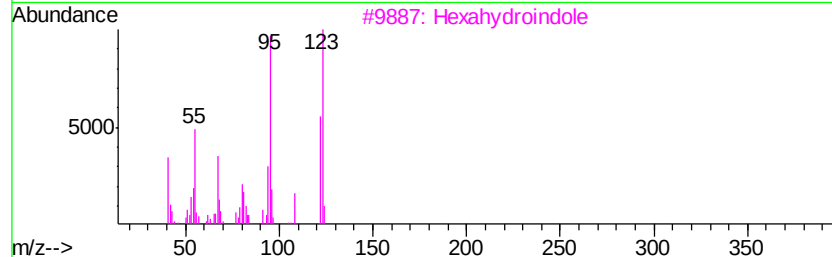
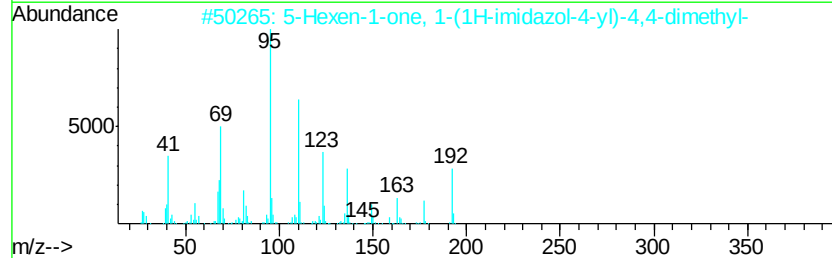
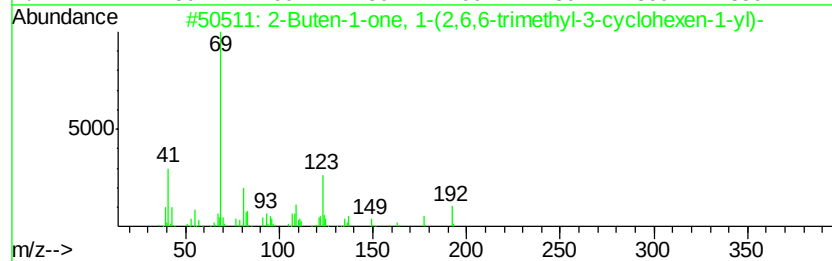
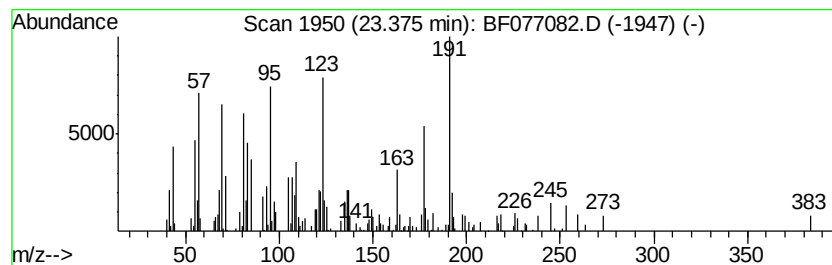
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 unknown23.37 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.37	16.32 ng	178315	Perylene-d12	22.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Buten-1-one, 1-(2,6,6-trimethy...	192	C13H20O	041436-42-4	35
2		5-Hexen-1-one, 1-(1H-imidazol-4-...	192	C11H16N2O	069393-41-5	30
3		Hexahydroindole	123	C8H13N	018159-32-5	27
4		2-Buten-1-one, 1-(2,6,6-trimethy...	192	C13H20O	024720-09-0	14
5		1-(3,3,6a-Trimethyl-1a,2,3,5,6a,...	220	C14H20O2	1000190-30-6	14



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012615\
 Data File : BF077082.D
 Acq On : 27 Jan 2015 9:59
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 Sample : G1160-03
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF011415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown3.83	3.83	12.6	ng	85293	1	7.76	134929	20.0
Benzophenone	16.47	10.2	ng	118454	4	17.64	233388	20.0
Methyl tetradecan...	18.46	176.3	ng	2057770	4	17.64	233388	20.0
9-Octadecenoic ac...	19.39	447.4	ng	5220900	4	17.64	233388	20.0
Octadecanoic acid...	19.50	46.1	ng	976629	5	21.15	423902	20.0
Ethanone, 1-cyclo...	20.31	4.4	ng	93898	5	21.15	423902	20.0
Eicosanoic acid, ...	20.43	13.6	ng	287649	5	21.15	423902	20.0
unknown20.49	20.49	5.6	ng	118909	5	21.15	423902	20.0
unknown20.68	20.68	7.3	ng	154021	5	21.15	423902	20.0
unknown20.95	20.95	6.5	ng	137376	5	21.15	423902	20.0
unknown22.04	22.04	12.6	ng	137369	6	22.78	218586	20.0
unknown23.37	23.37	16.3	ng	178315	6	22.78	218586	20.0