

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062016\
 Data File : BF088214.D
 Acq On : 21 Jun 2016 12:20
 Operator : UM/SJ
 Sample : H3580-29
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P8-B4(0-12)C

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-BF060716.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	1.656	5	6	14	rVV	175014	272336	9.29%	1.347%
2	1.770	14	16	54	rVB	1418685	2932923	100.00%	14.505%
3	4.822	281	283	294	rBB	69273	112911	3.85%	0.558%
4	5.267	319	322	324	rBV	1336929	1715149	58.48%	8.482%
5	6.330	412	415	417	rBV	1309126	1779920	60.69%	8.803%
6	6.445	422	425	427	rBV	1353580	1770622	60.37%	8.757%
7	6.662	443	444	447	rBV	368059	456921	15.58%	2.260%
8	6.822	456	458	461	rBV	1581298	1371965	46.78%	6.785%
9	7.233	492	494	497	rBV	866831	981499	33.46%	4.854%
10	7.953	555	557	559	rBV	749515	625680	21.33%	3.094%
11	9.028	649	651	654	rBV	1868338	1845772	62.93%	9.128%
12	9.428	684	686	688	rBV	567916	458340	15.63%	2.267%
13	9.645	703	705	707	rBV	62818	49674	1.69%	0.246%
14	9.702	707	710	712	rBV	692655	634501	21.63%	3.138%
15	10.136	747	748	750	rVB	74391	78063	2.66%	0.386%
16	10.354	764	767	771	rBV	28083	56214	1.92%	0.278%
17	10.491	776	779	781	rBV	759910	827187	28.20%	4.091%
18	10.616	788	790	794	rBV	75034	126375	4.31%	0.625%
19	11.062	827	829	830	rBV	38777	51779	1.77%	0.256%
20	11.176	837	839	841	rBV	610246	545389	18.60%	2.697%
21	11.782	890	892	897	rVB3	14877	31379	1.07%	0.155%
22	12.171	923	926	928	rBV2	15441	32446	1.11%	0.160%
23	12.274	933	935	937	rVB2	39312	43180	1.47%	0.214%
24	12.388	943	945	946	rBV	63769	60864	2.08%	0.301%
25	12.617	961	965	966	rBV2	45171	89685	3.06%	0.444%
26	12.754	975	977	980	rBV	1068729	1311110	44.70%	6.484%
27	12.994	996	998	1000	rBV2	46440	65189	2.22%	0.322%
28	13.040	1000	1002	1005	rVV3	21578	38347	1.31%	0.190%
29	13.131	1007	1010	1012	rBV2	22810	50110	1.71%	0.248%
30	13.200	1014	1016	1018	rBV3	18785	30326	1.03%	0.150%
31	13.245	1018	1020	1022	rBV2	19567	31677	1.08%	0.157%
32	13.291	1022	1024	1026	rBV2	25829	35099	1.20%	0.174%
33	13.337	1026	1028	1031	rBV	58073	76681	2.61%	0.379%
34	13.668	1055	1057	1058	rBV	60305	61927	2.11%	0.306%

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P8-B4(0-12)C

Integration Parameters: rteint.p
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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 >
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35	13.805	1067	1069	1075	rBV	552119	587499	20.03%	2.906%
36	13.977	1082	1084	1089	rBV	64967	109018	3.72%	0.539%
37	14.285	1109	1111	1112	rBV	65300	63847	2.18%	0.316%
38	14.583	1135	1137	1138	rBV	82655	117554	4.01%	0.581%
39	15.188	1188	1190	1195	rVB	285033	474215	16.17%	2.345%
40	15.577	1222	1224	1226	rBV	64896	101043	3.45%	0.500%
41	16.137	1271	1273	1279	rVB2	59740	115602	3.94%	0.572%

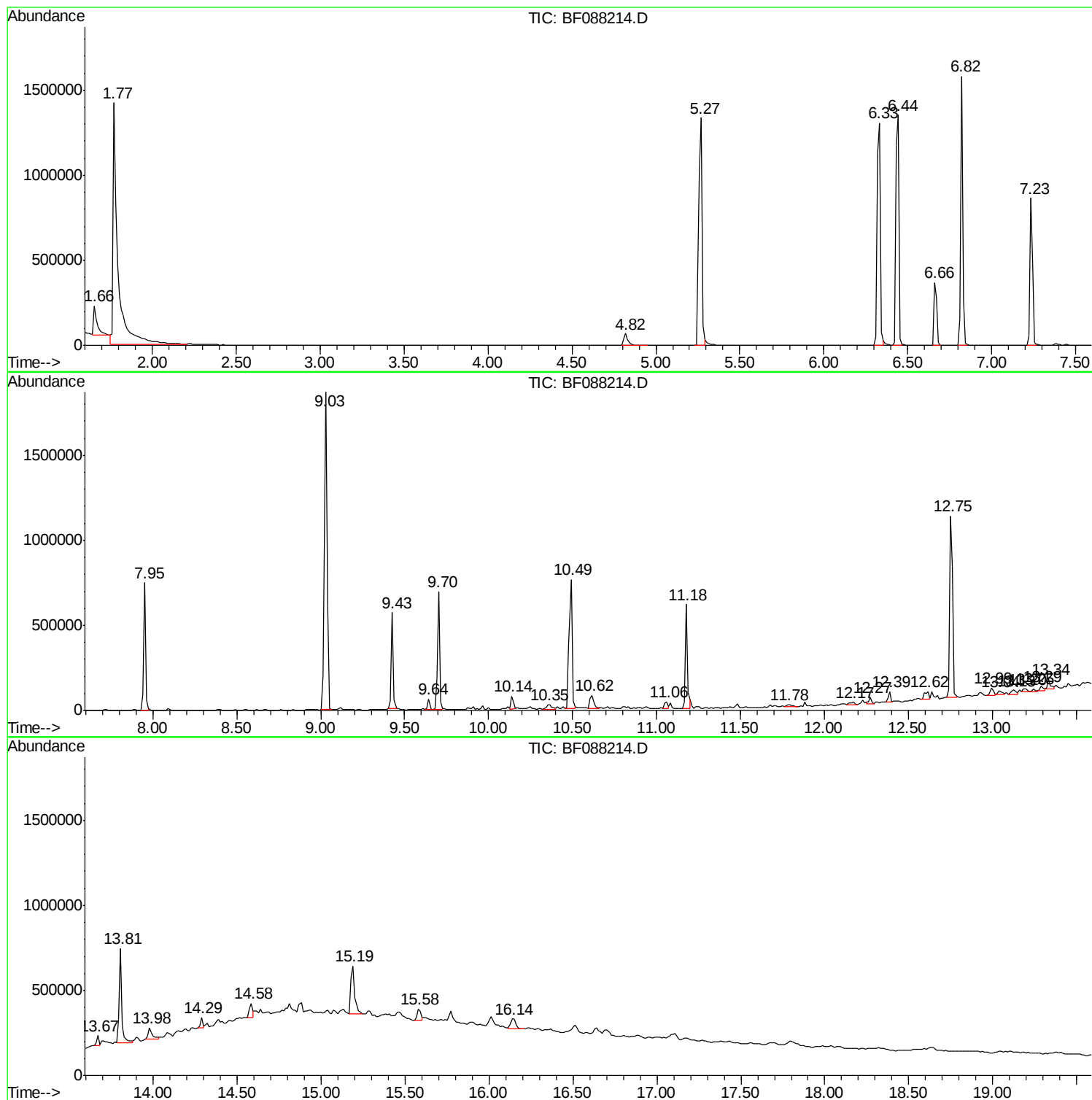
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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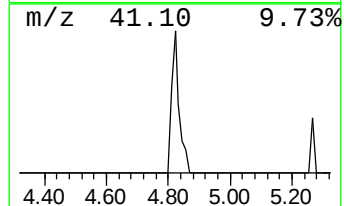
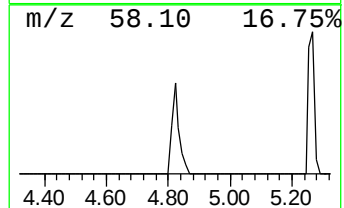
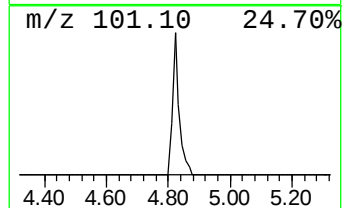
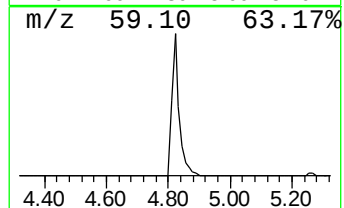
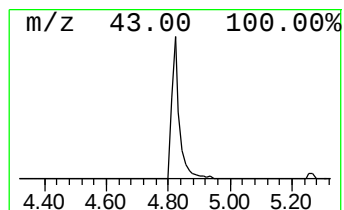
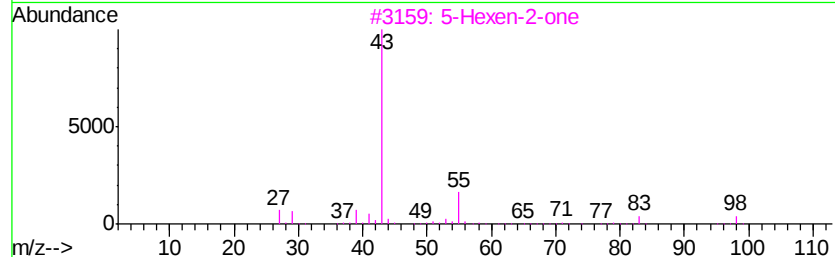
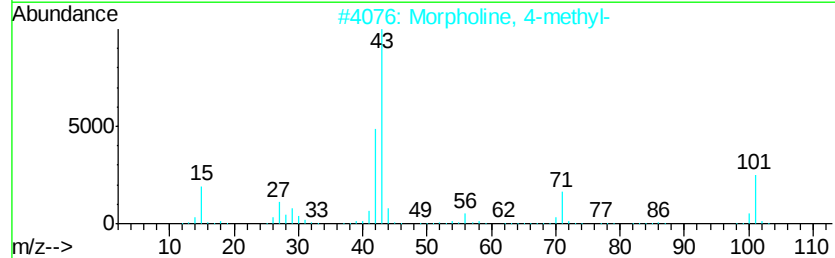
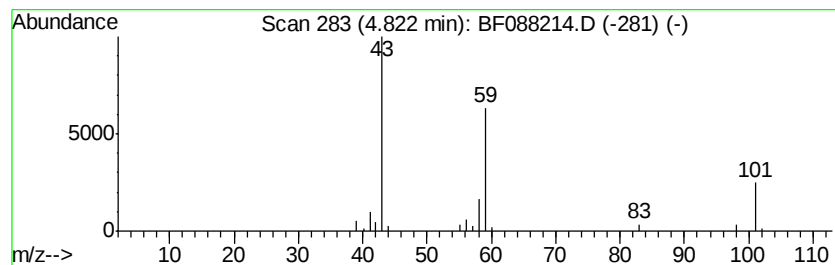
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Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.82	4.94 ng	112911	1,4-Dichlorobenzene-d4	6.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
3		5-Hexen-2-one	98	C6H10O	000109-49-9	9
4		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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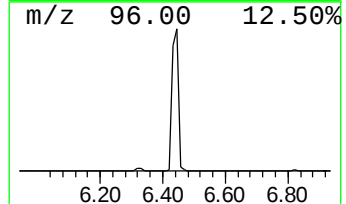
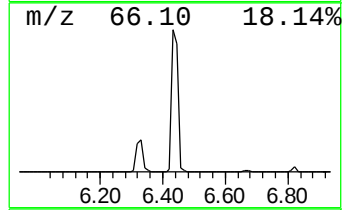
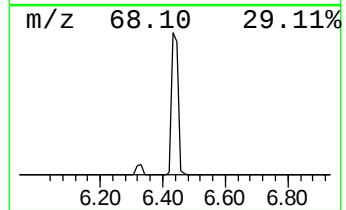
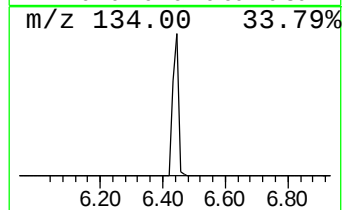
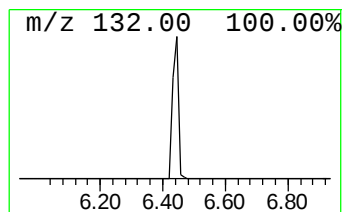
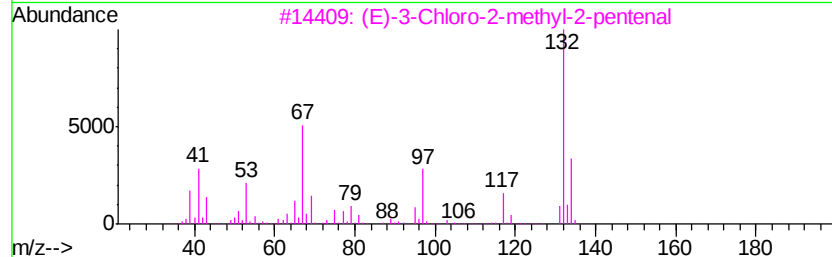
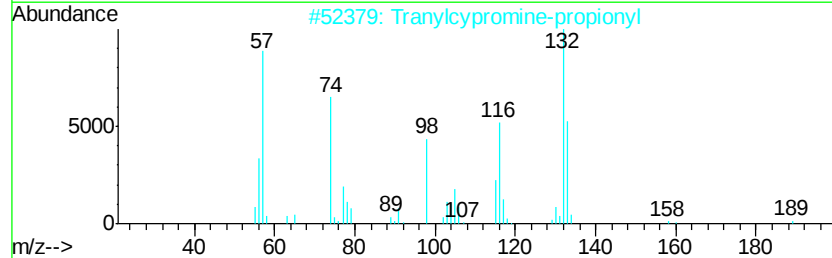
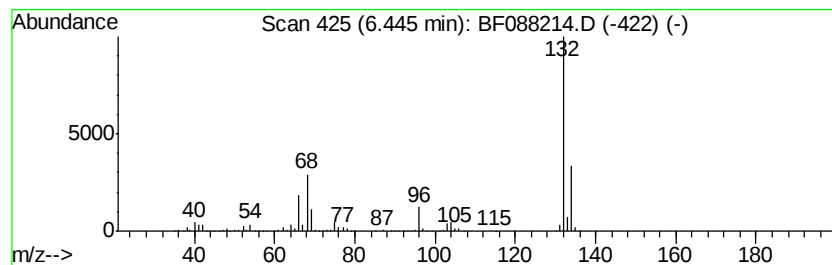
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
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TIC Library : C:\Database\NIST11.L
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Peak Number 4 3-Chloro-6-fluoro-pyrazine Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.44	77.50 ng	1770620	1,4-Dichlorobenzene-d4	6.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	50
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	37
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	33
4		Amphetaminil	250	C17H18N2	017590-01-1	28
5		1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	25



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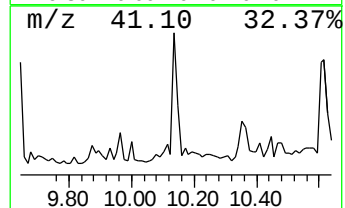
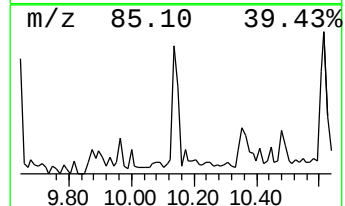
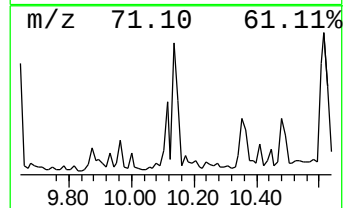
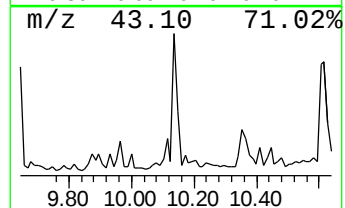
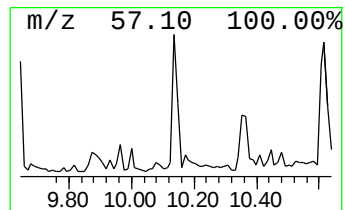
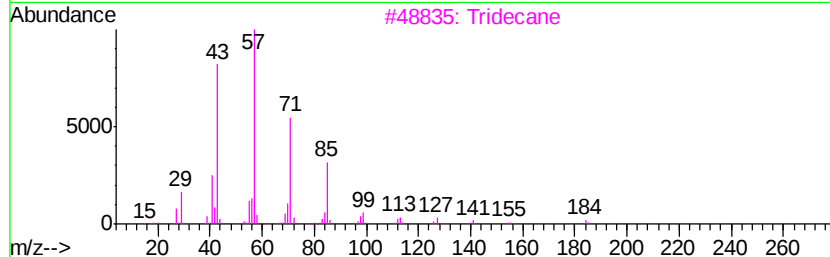
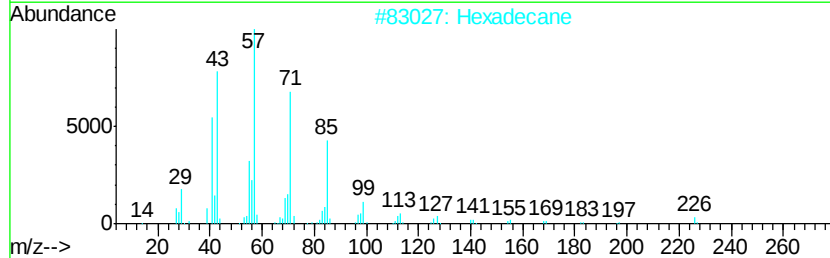
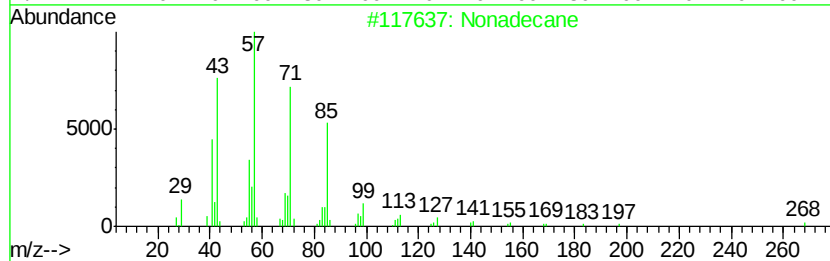
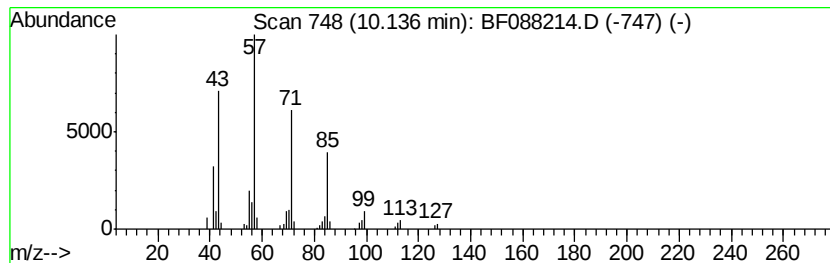
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TIC Library : C:\Database\NIST11.L
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Peak Number 6 Nonadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.14	2.46 ng	78063	Acenaphthene-d10	9.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	90
2		Hexadecane	226	C16H34	000544-76-3	87
3		Tridecane	184	C13H28	000629-50-5	86
4		Tetradecane	198	C14H30	000629-59-4	86
5		Pentadecane	212	C15H32	000629-62-9	78



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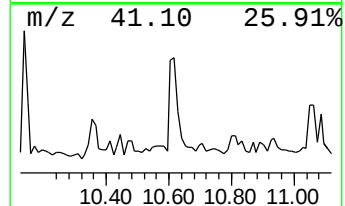
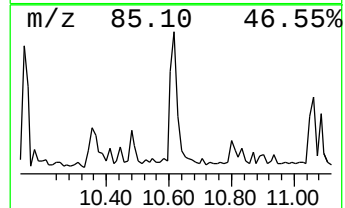
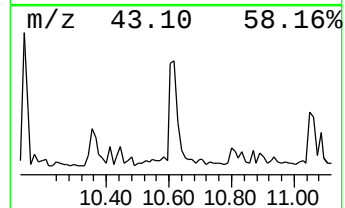
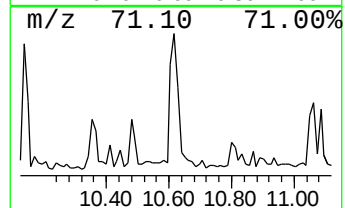
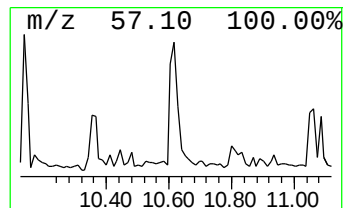
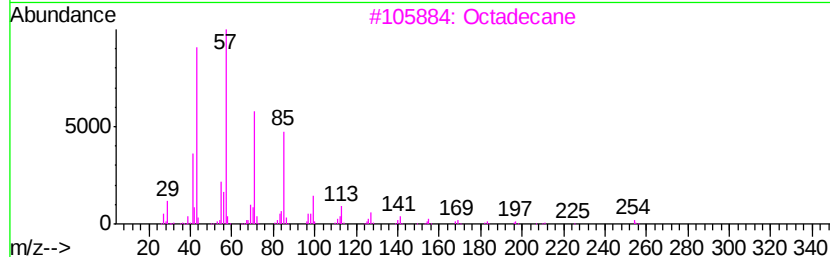
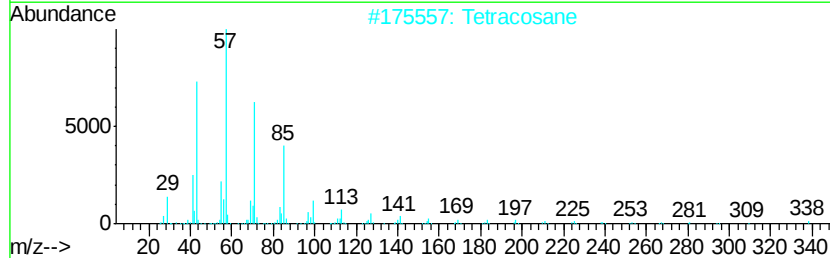
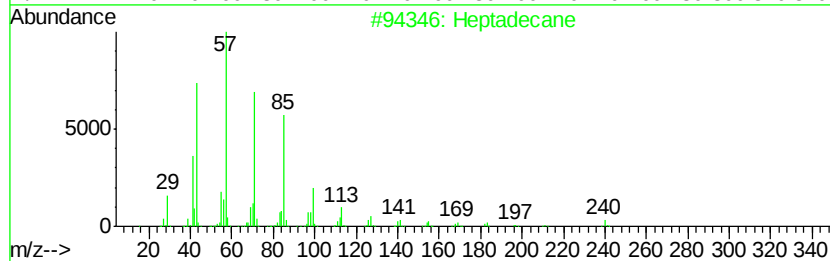
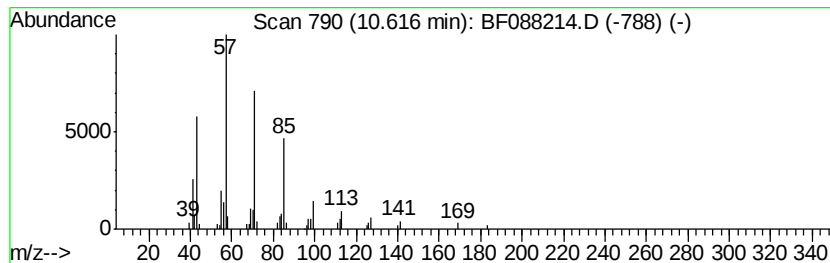
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TIC Library : C:\Database\NIST11.L
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Peak Number 7 Heptadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.62	4.63 ng	126375	Phenanthrene-d10	11.18

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	91
2	Tetracosane		338	C24H50	000646-31-1	91
3	Octadecane		254	C18H38	000593-45-3	91
4	Heptadecane, 2,6,10,15-tetramethyl-		296	C21H44	054833-48-6	91
5	Hexadecane		226	C16H34	000544-76-3	91



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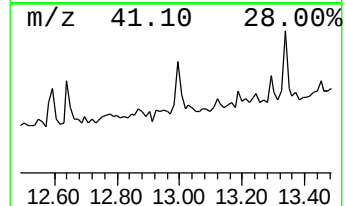
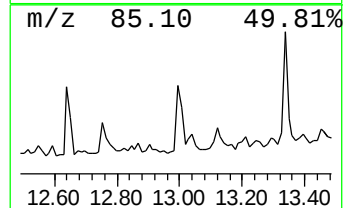
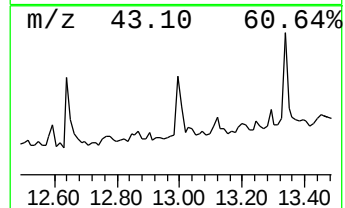
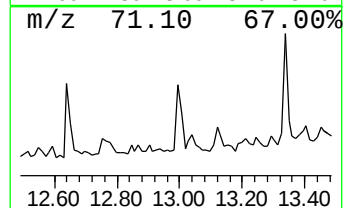
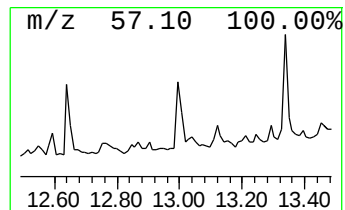
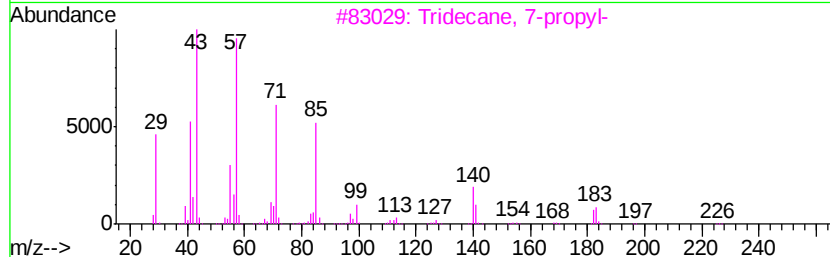
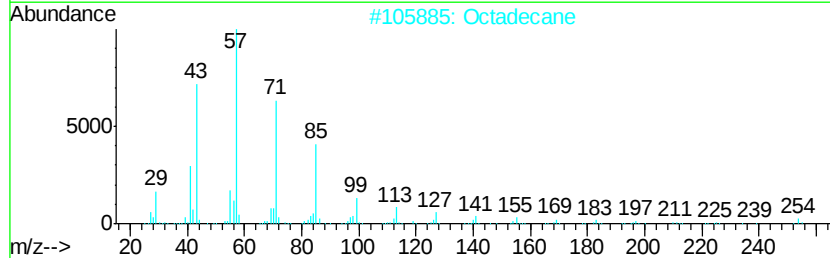
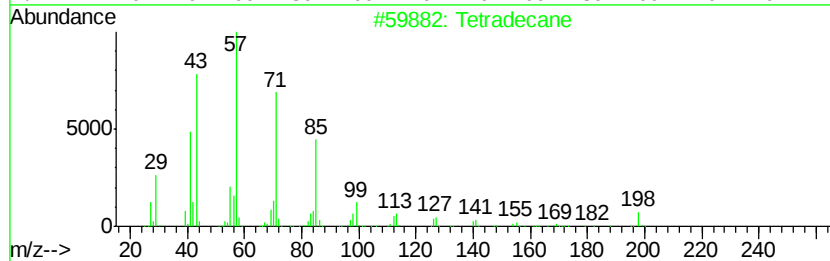
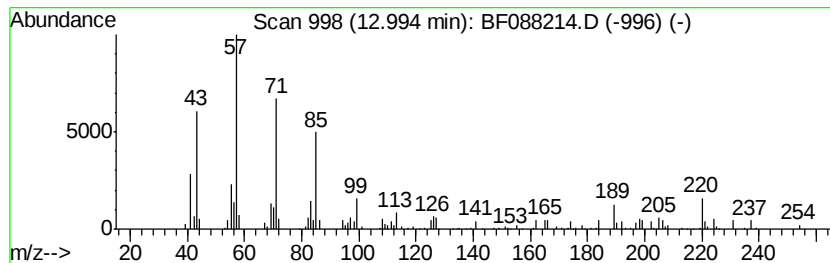
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TIC Library : C:\Database\NIST11.L
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Peak Number 9 Tetradecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.99	2.22 ng	65189	Chrysene-d12	13.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane	198	C14H30	000629-59-4	83
2		Octadecane	254	C18H38	000593-45-3	78
3		Tridecane, 7-propyl-	226	C16H34	055045-09-5	78
4		Hexadecane, 7,9-dimethyl-	254	C18H38	021164-95-4	78
5		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	60



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P8-B4(0-12)C

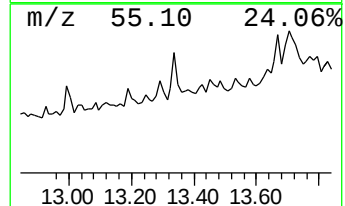
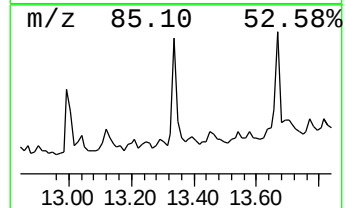
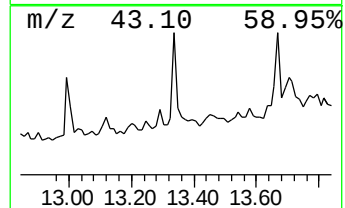
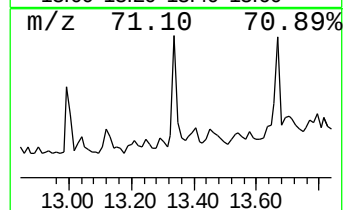
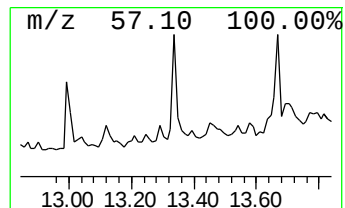
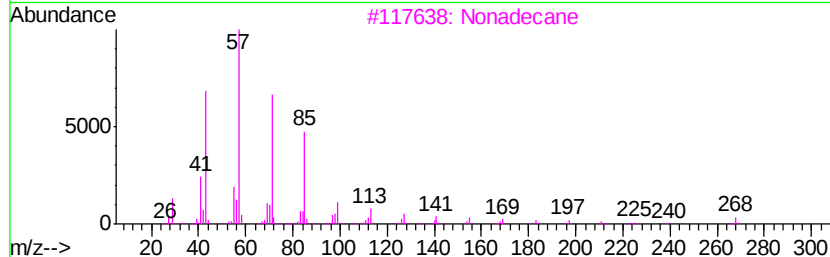
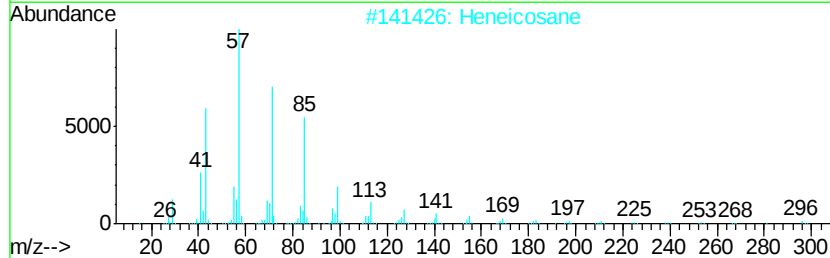
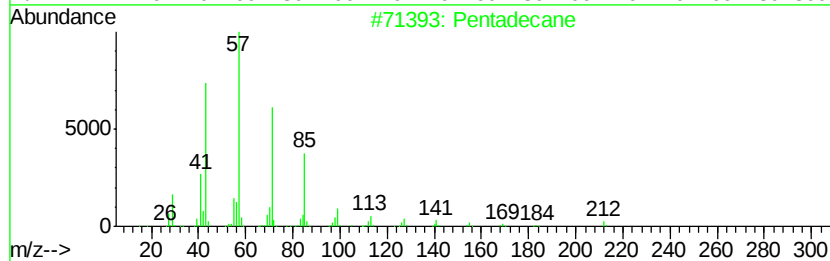
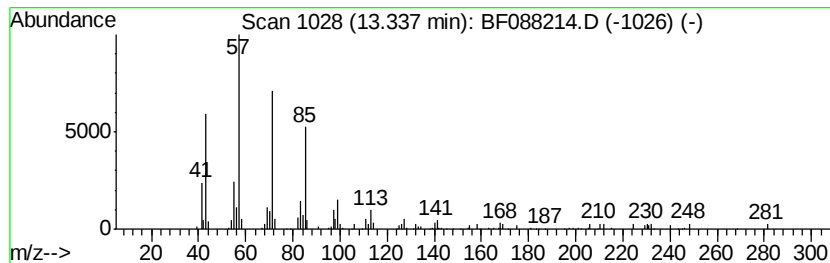
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Pentadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.34	2.61 ng	76681	Chrysene-d12	13.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane	212	C15H32	000629-62-9	91
2			Heneicosane	296	C21H44	000629-94-7	86
3			Nonadecane	268	C19H40	000629-92-5	86
4			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	80
5			Octadecane	254	C18H38	000593-45-3	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062016\
Data File : BF088214.D
Acq On : 21 Jun 2016 12:20
Operator : UM/SJ
Sample : H3580-29
Misc :
ALS Vial : 43 Sample Multiplier: 1

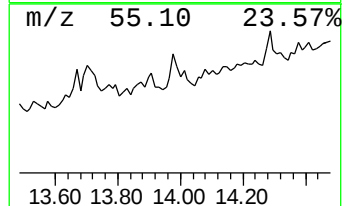
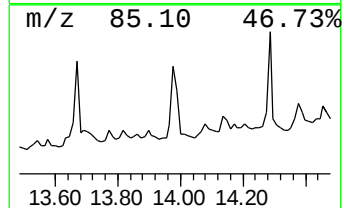
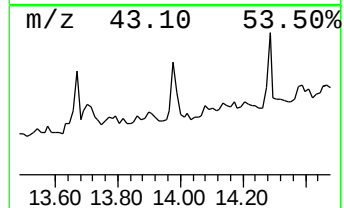
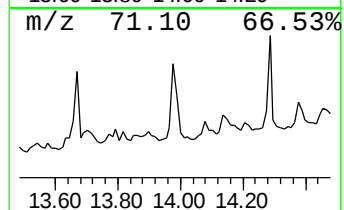
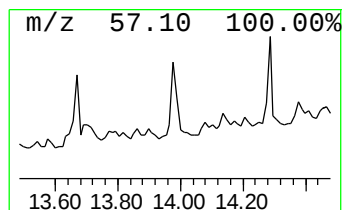
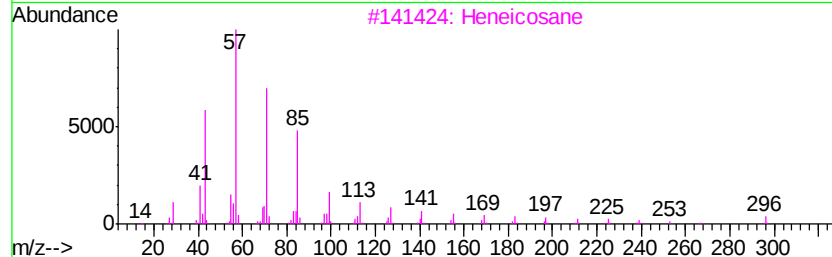
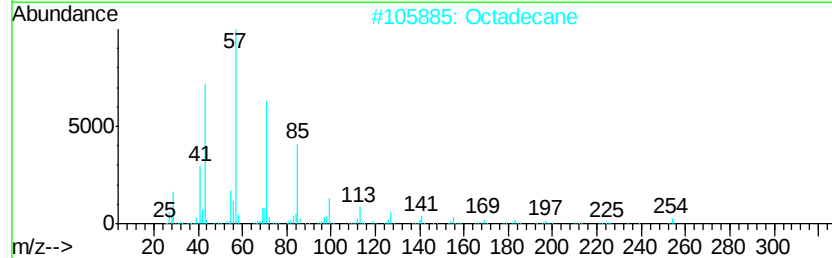
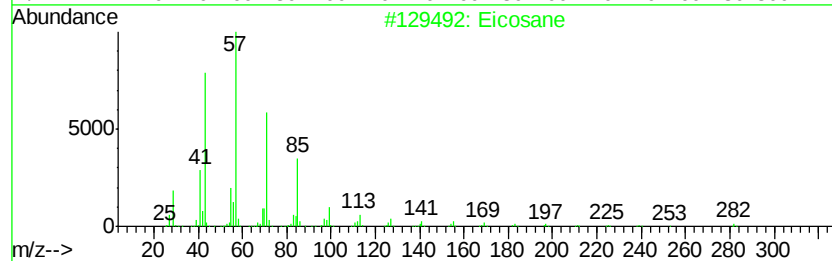
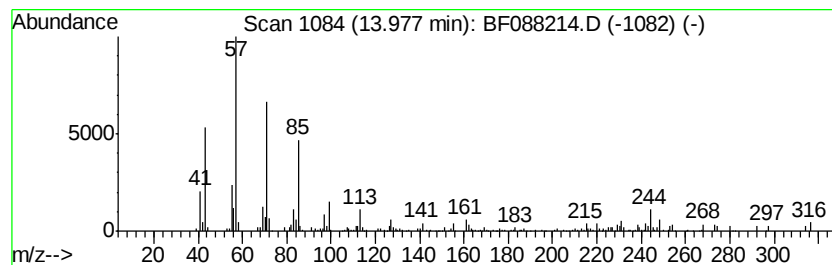
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Eicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.98	3.71 ng	109018	Chrysene-d12		13.81
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	96
2	Octadecane	254	C18H38	000593-45-3	92
3	Heneicosane	296	C21H44	000629-94-7	90
4	Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	81
5	Octacosane	394	C28H58	000630-02-4	74



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062016\
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Operator : UM/SJ
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Instrument :
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ClientSampleId :
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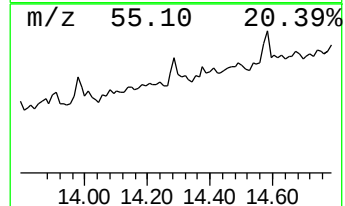
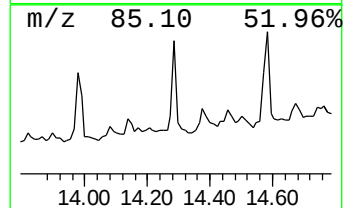
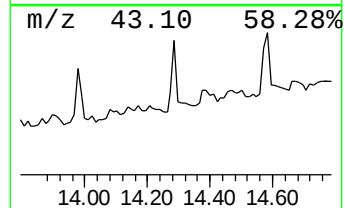
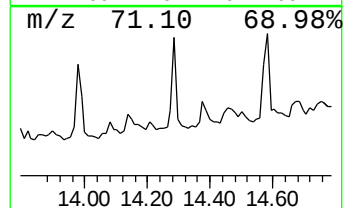
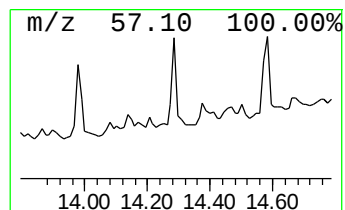
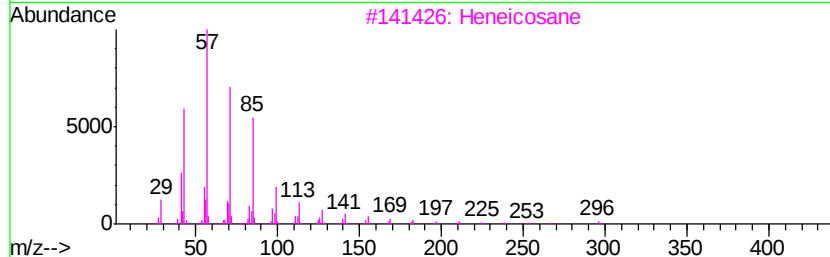
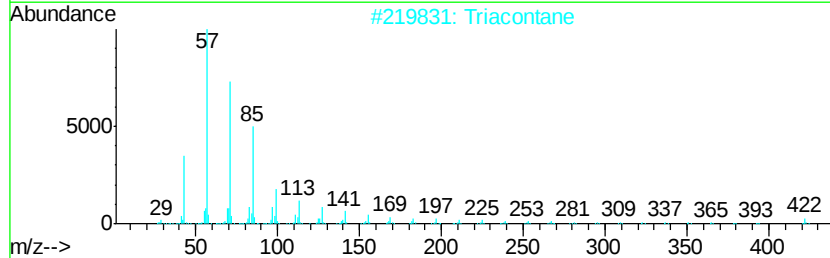
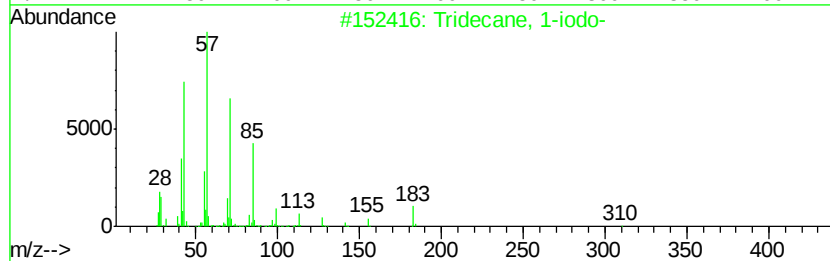
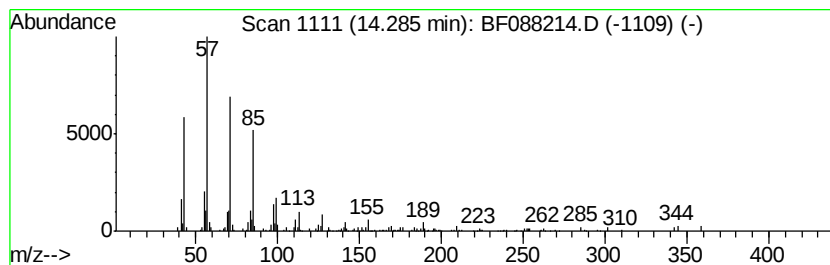
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Tridecane, 1-iodo- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.29	2.17 ng	63847	Chrysene-d12	13.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	91
2		Triacontane	422	C30H62	000638-68-6	87
3		Heneicosane	296	C21H44	000629-94-7	87
4		Docosane	310	C22H46	000629-97-0	87
5		Pentacosane	352	C25H52	000629-99-2	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062016\
Data File : BF088214.D
Acq On : 21 Jun 2016 12:20
Operator : UM/SJ
Sample : H3580-29
Misc :
ALS Vial : 43 Sample Multiplier: 1

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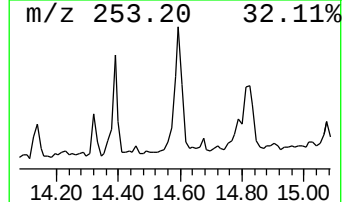
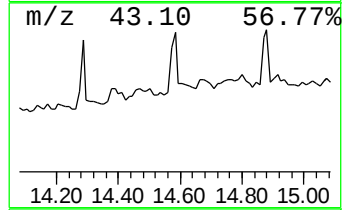
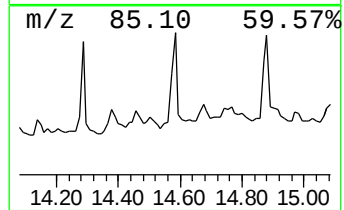
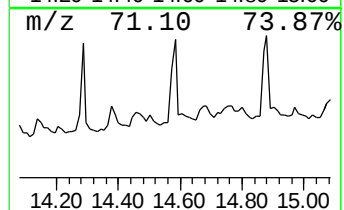
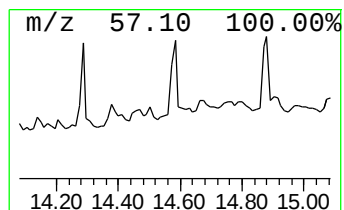
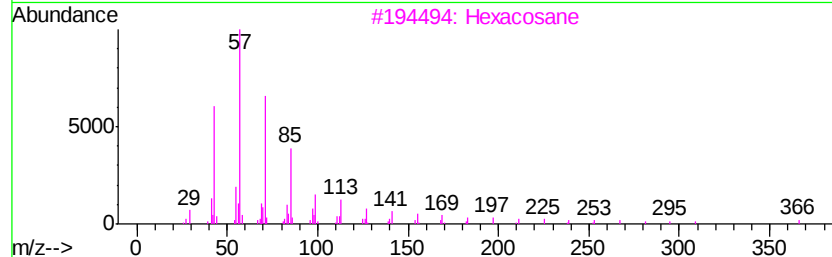
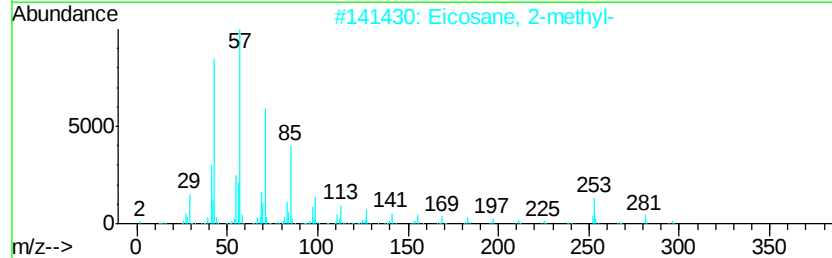
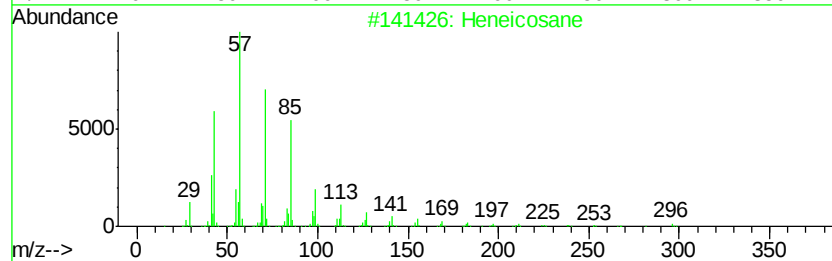
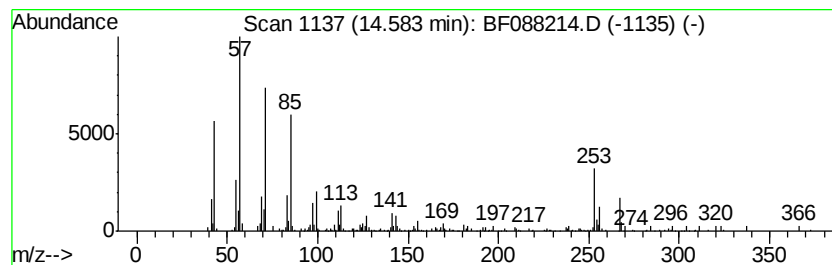
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Heneicosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.58	4.96 ng	117554	Perylene-d12	15.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	89
2		Eicosane, 2-methyl-	296	C21H44	001560-84-5	81
3		Hexacosane	366	C26H54	000630-01-3	78
4		Octadecane	254	C18H38	000593-45-3	76
5		2-methyloctacosane	408	C29H60	1000376-72-8	76



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062016\
Data File : BF088214.D
Acq On : 21 Jun 2016 12:20
Operator : UM/SJ
Sample : H3580-29
Misc :
ALS Vial : 43 Sample Multiplier: 1

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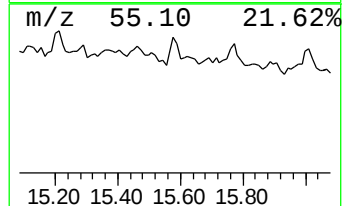
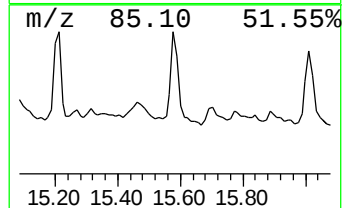
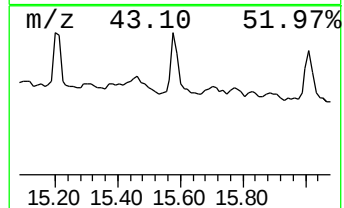
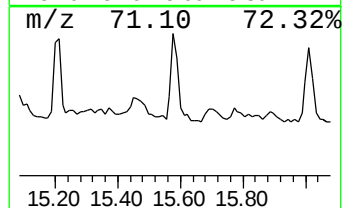
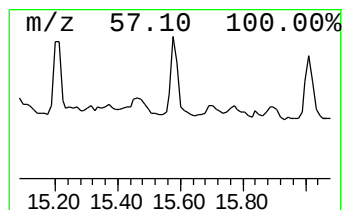
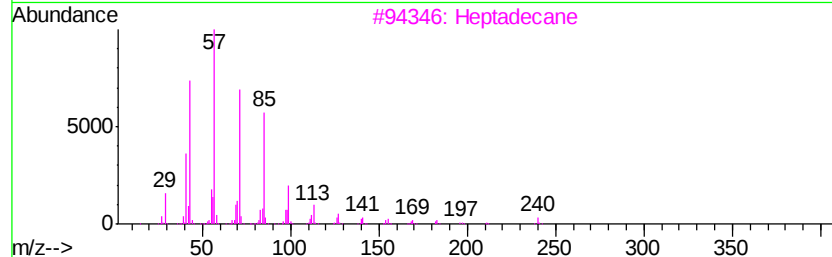
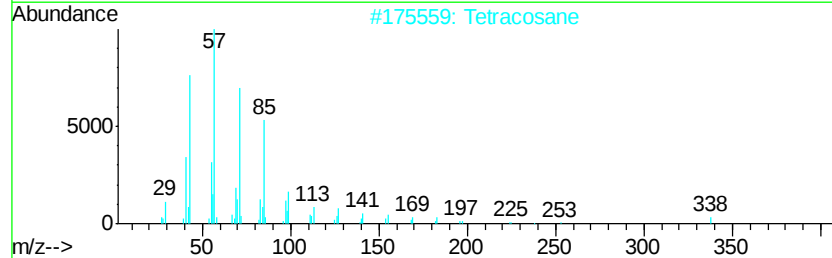
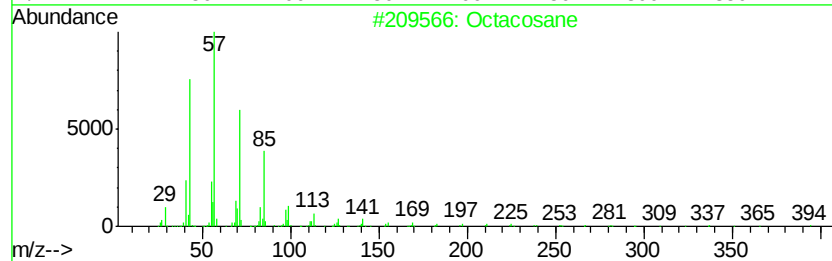
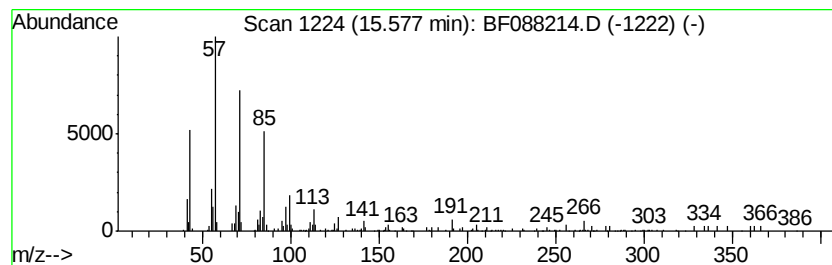
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Octacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.58	4.26 ng	101043	Perylene-d12	15.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	90
2		Tetracosane	338	C24H50	000646-31-1	87
3		Heptadecane	240	C17H36	000629-78-7	87
4		Heneicosane	296	C21H44	000629-94-7	87
5		Hexacosane	366	C26H54	000630-01-3	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062016\
Data File : BF088214.D
Acq On : 21 Jun 2016 12:20
Operator : UM/SJ
Sample : H3580-29
Misc :
ALS Vial : 43 Sample Multiplier: 1

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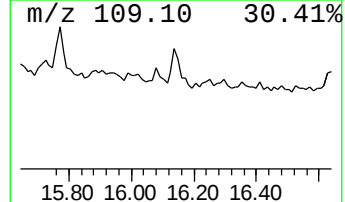
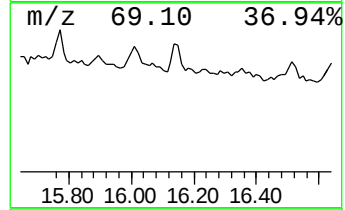
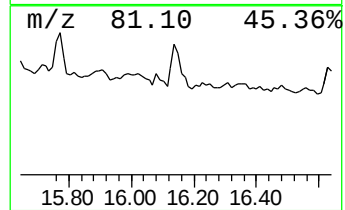
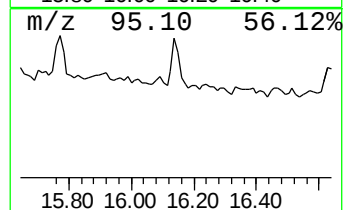
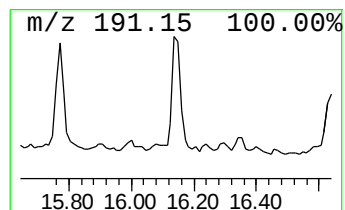
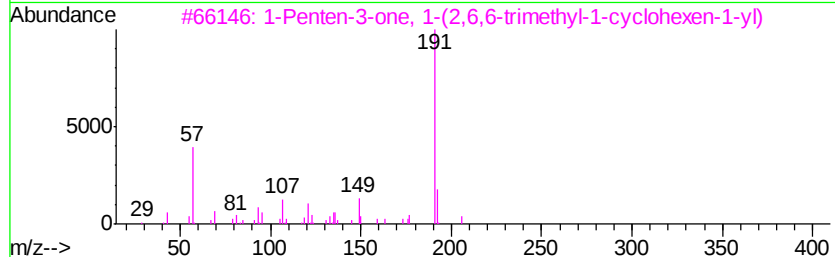
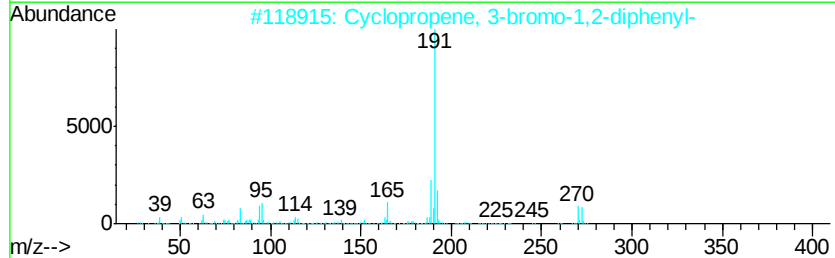
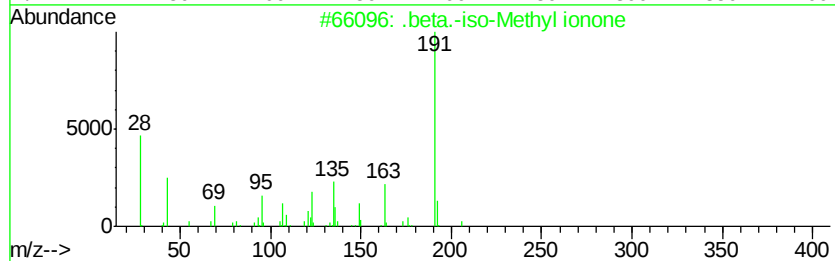
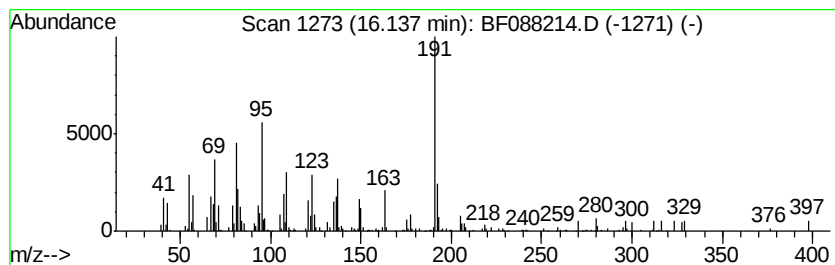
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 .beta.-iso-Methyl ionone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.14	4.88 ng	115602	Perylene-d12	15.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	76
2		Cyclopropene, 3-bromo-1,2-diphenyl-	270	C15H11Br	029559-17-9	38
3		1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	38
4		Benzene, 1,1'-[1,2-cyclohexanedi...	300	C18H20S2	1000362-19-6	38
5		2,4,5,5,8a-Pentamethyl-6,7,8,8a-...	206	C14H22O	1000195-40-9	37



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062016\
Data File : BF088214.D
Acq On : 21 Jun 2016 12:20
Operator : UM/SJ
Sample : H3580-29
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.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
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3-Chloro-6-fluoro...	6.44	77.5	ng	1770620	1	6.66	456921	20.0
Nonadecane	10.14	2.5	ng	78063	3	9.70	634501	20.0
Heptadecane	10.62	4.6	ng	126375	4	11.18	545389	20.0
Tetradecane	12.99	2.2	ng	65189	5	13.81	587499	20.0
Pentadecane	13.34	2.6	ng	76681	5	13.81	587499	20.0
Eicosane	13.98	3.7	ng	109018	5	13.81	587499	20.0
Tridecane, 1-iodo-	14.29	2.2	ng	63847	5	13.81	587499	20.0
Heneicosane	14.58	5.0	ng	117554	6	15.19	474215	20.0
Octacosane	15.58	4.3	ng	101043	6	15.19	474215	20.0
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