

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG071719\
 Data File : BG041699.D
 Acq On : 17 Jul 2019 22:17
 Operator : HP/JU
 Sample : K3835-08
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-4(7-9)

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_G\METHODS\8270-BG071519.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	3.199	13	19	34	rVB	363625	674751	10.09%	1.661%
2	5.608	422	429	440	rBV	1866391	2991792	44.75%	7.363%
3	7.200	692	700	717	rBV	1913504	3320773	49.67%	8.173%
4	7.576	756	764	773	rBV	1844414	3207988	47.98%	7.895%
5	8.046	837	844	851	rBV	378953	634597	9.49%	1.562%
6	8.369	891	899	912	rBV	1398179	2431331	36.37%	5.984%
7	9.204	1033	1041	1053	rBV	1084356	1997223	29.87%	4.916%
8	10.855	1313	1322	1331	rVB	469370	868305	12.99%	2.137%
9	13.293	1730	1737	1751	rBV	2735424	4221895	63.15%	10.391%
10	14.110	1870	1876	1887	rVB	308912	472180	7.06%	1.162%
11	14.662	1962	1970	1981	rVB2	705988	1171574	17.52%	2.883%
12	16.143	2214	2222	2236	rBV	2436469	3821198	57.15%	9.405%
13	17.400	2429	2436	2444	rBV2	986925	1581167	23.65%	3.892%
14	20.003	2873	2879	2892	rBV	4786525	6685883	100.00%	16.455%
15	21.325	3098	3104	3130	rBV3	130652	562567	8.41%	1.385%
16	21.648	3152	3159	3169	rBV2	1186362	1969473	29.46%	4.847%
17	21.871	3192	3197	3205	rBV	68769	105146	1.57%	0.259%
18	22.482	3296	3301	3309	rBV	124089	203320	3.04%	0.500%
19	23.181	3413	3420	3426	rVB	167363	329582	4.93%	0.811%
20	23.992	3551	3558	3570	rVB	184123	413433	6.18%	1.018%
21	24.844	3693	3703	3713	rBV2	708943	2005116	29.99%	4.935%
22	24.950	3713	3721	3734	rVB	166549	420899	6.30%	1.036%
23	26.090	3906	3915	3926	rVB2	117057	353738	5.29%	0.871%
24	27.465	4142	4149	4159	rVB3	55644	186728	2.79%	0.460%

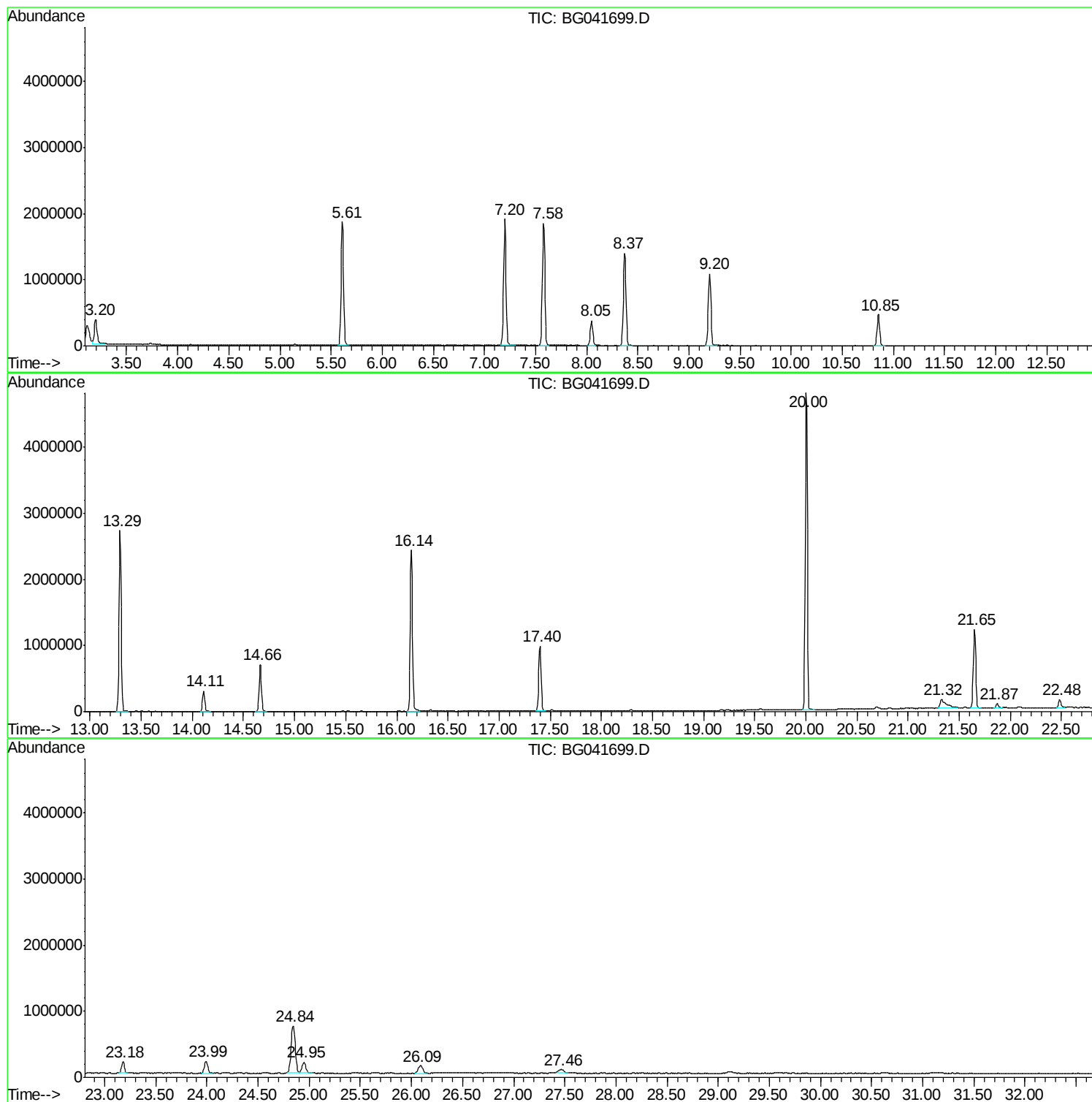
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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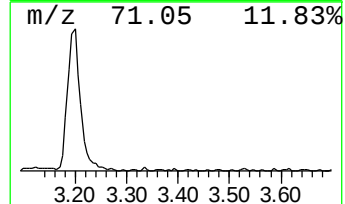
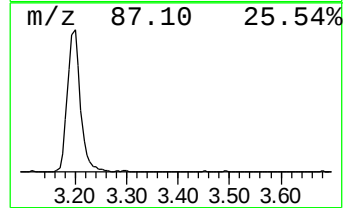
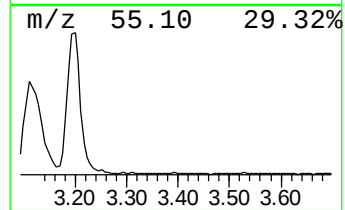
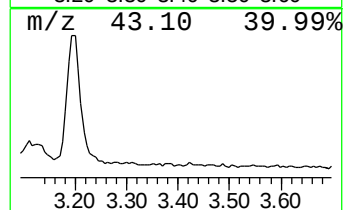
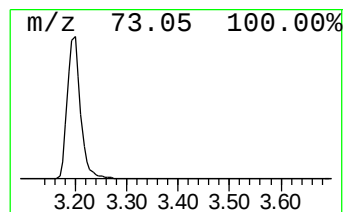
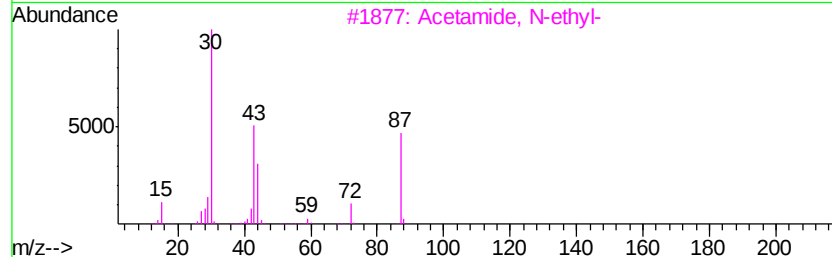
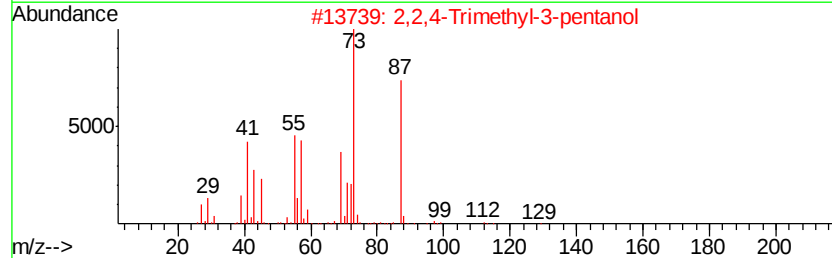
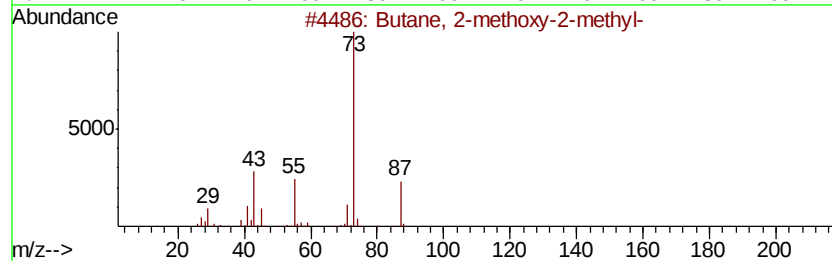
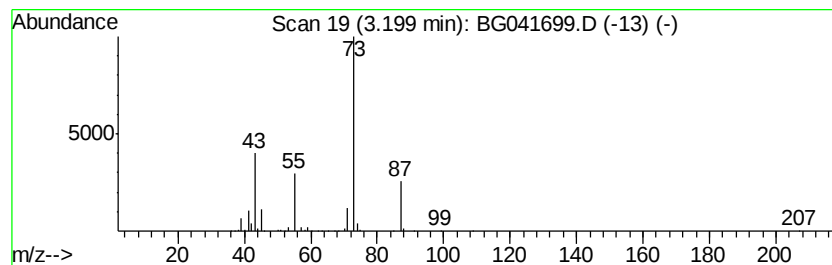
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.20	21.27 ng	674751	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	27
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
5		3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	23



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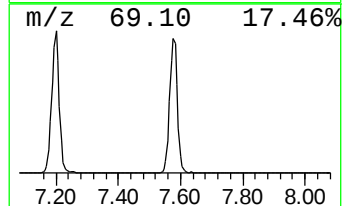
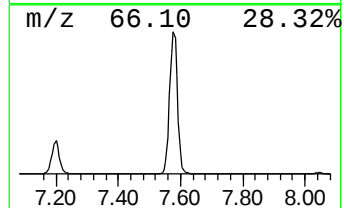
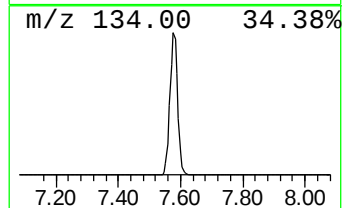
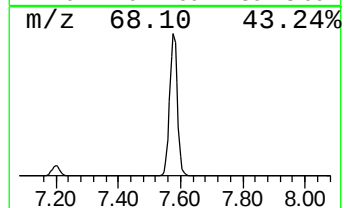
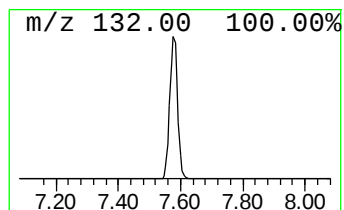
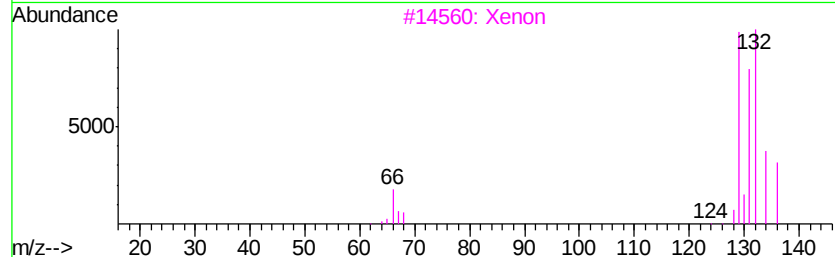
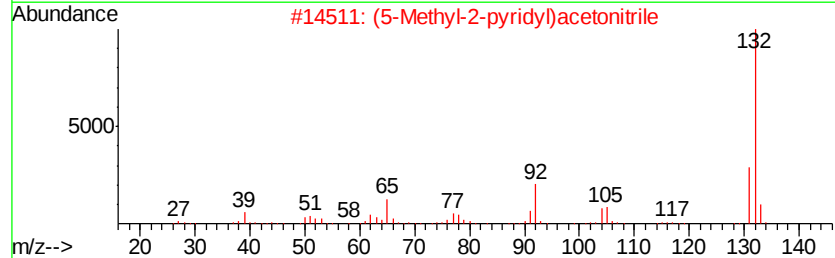
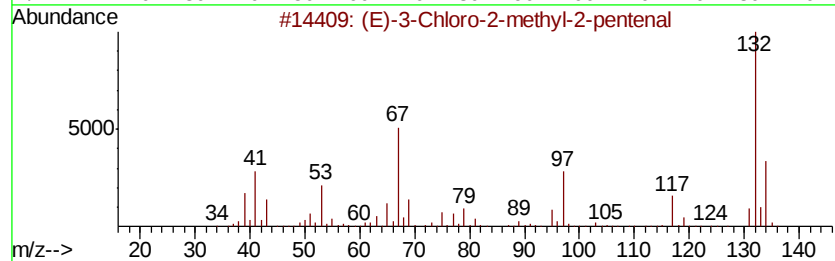
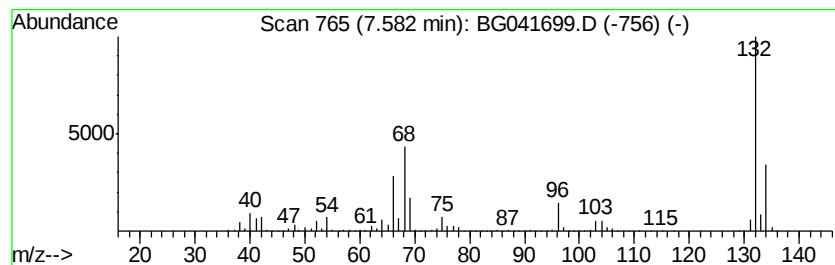
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	101.10 ng	3207990	1,4-Dichlorobenzene-d4	8.05

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
2			(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
3			Xenon	132	Xe	007440-63-3	9
4			1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5			1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



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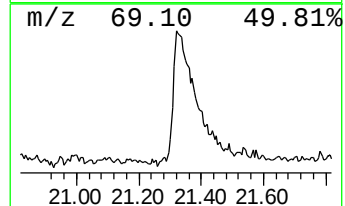
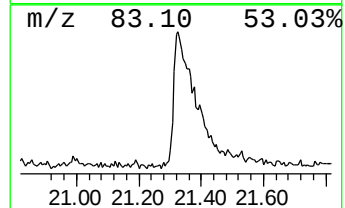
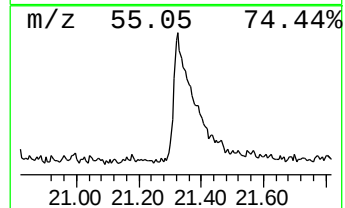
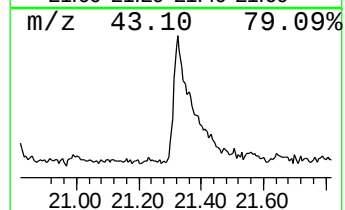
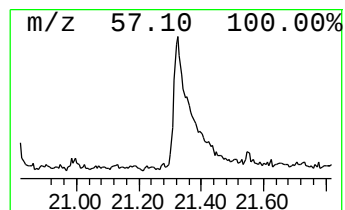
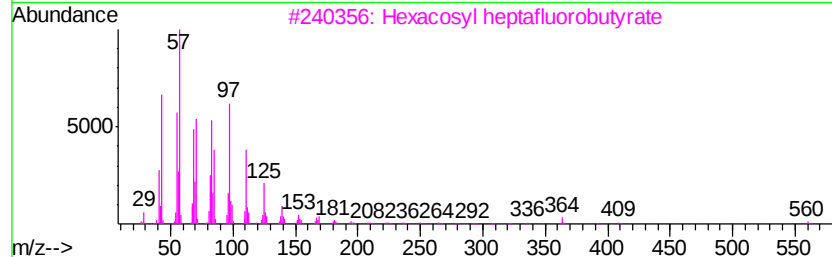
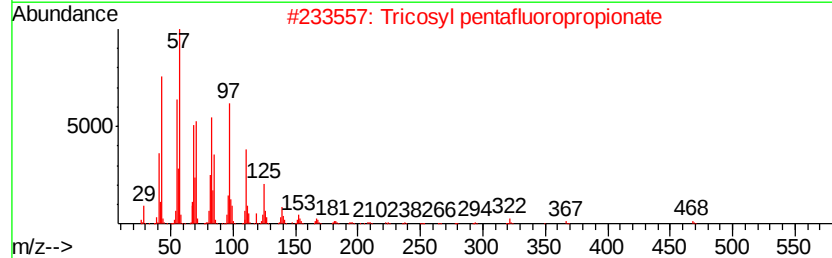
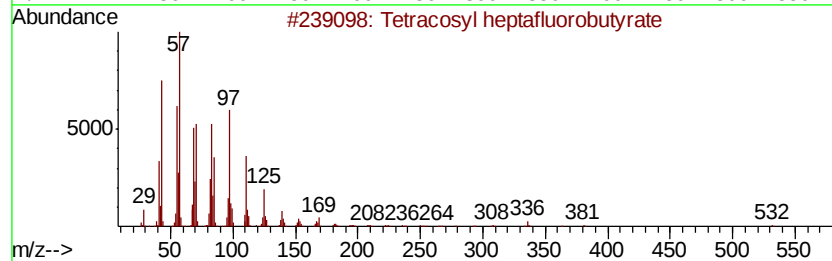
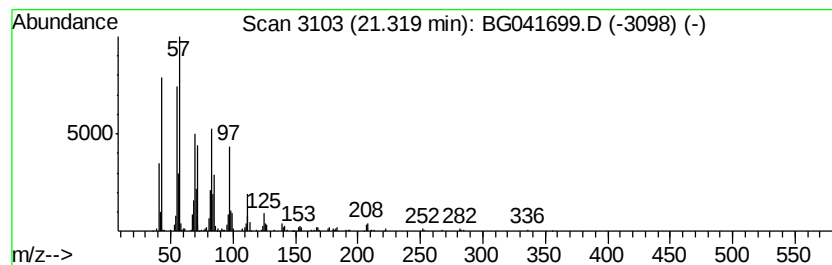
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Peak Number 4 Tetracosyl heptafluorobutyrate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.		
21.32	5.71 ng	562567	Chrysene-d12		21.65		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosyl heptafluorobutvrate	550	C28H49F7O2	1000351-83-7	91
2			Tricosyl pentafluoropropionate	486	C26H47F5O2	1000351-81-0	91
3			Hexacosyl heptafluorobutvrate	578	C30H53F7O2	1000351-83-3	91
4			Dotriacontyl heptafluorobutvrate	662	C36H65F7O2	1000351-84-2	91
5			Heptacosyl pentafluoropropionate	542	C30H55F5O2	1000351-81-6	91



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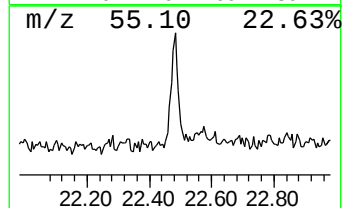
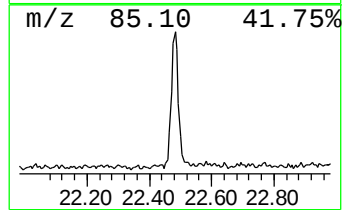
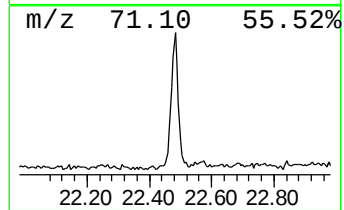
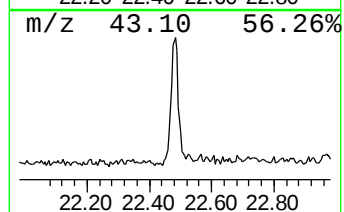
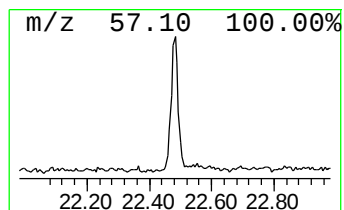
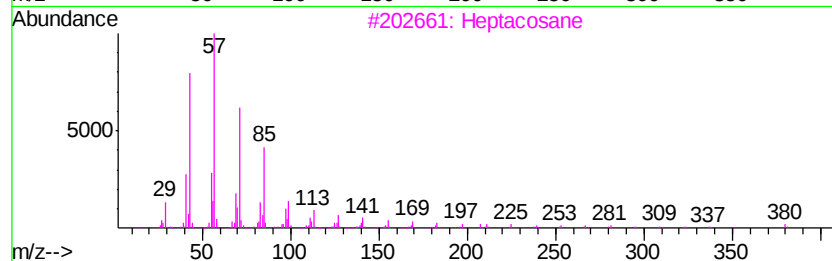
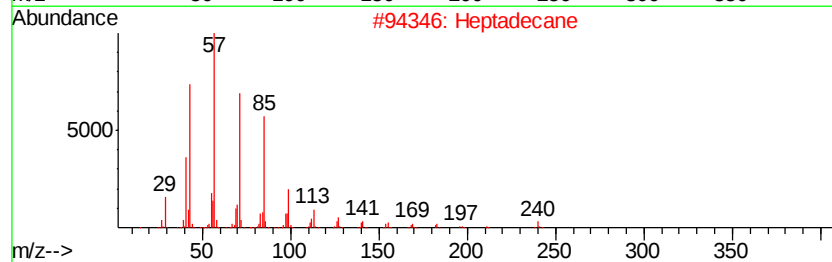
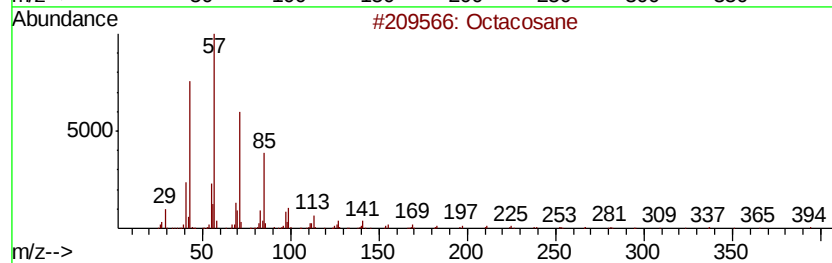
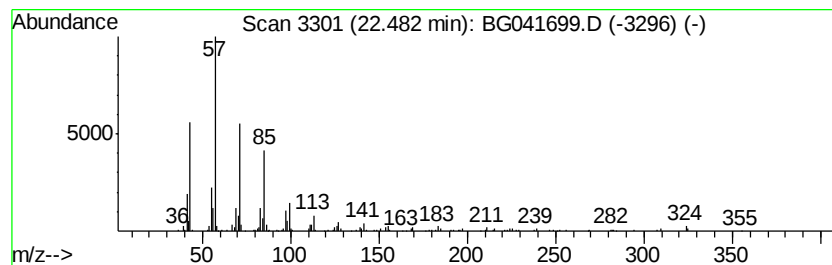
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Peak Number 5 Octacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.48	2.06 ng	203320	Chrysene-d12	21.65

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane		394	C28H58	000630-02-4	90
2	Heptadecane		240	C17H36	000629-78-7	90
3	Heptacosane		380	C27H56	000593-49-7	87
4	Tetracosane		338	C24H50	000646-31-1	87
5	Pentacosane		352	C25H52	000629-99-2	87



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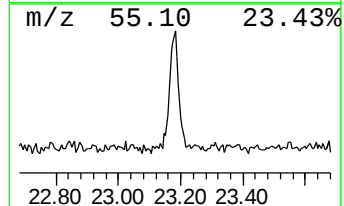
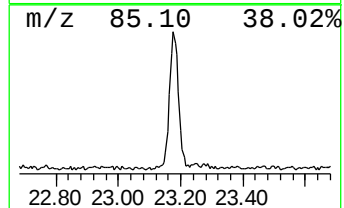
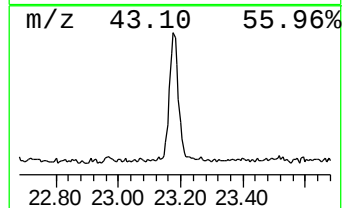
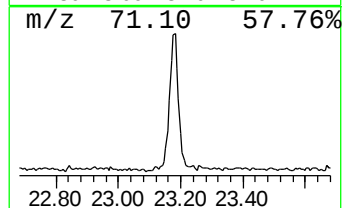
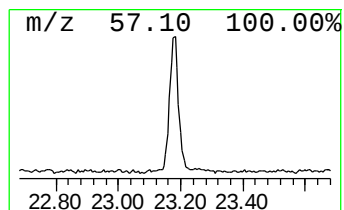
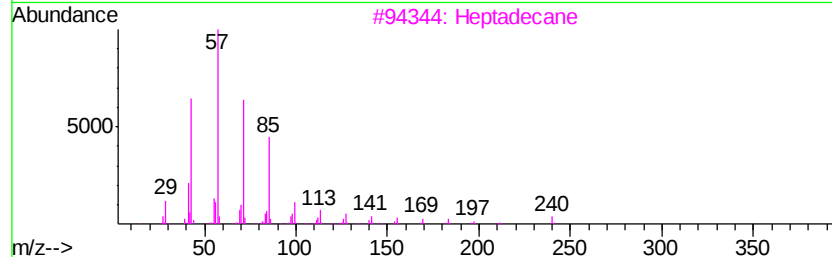
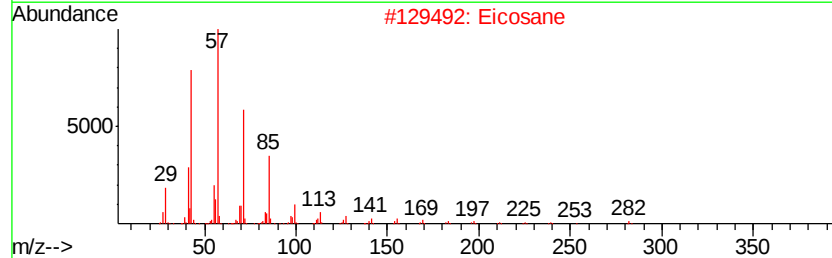
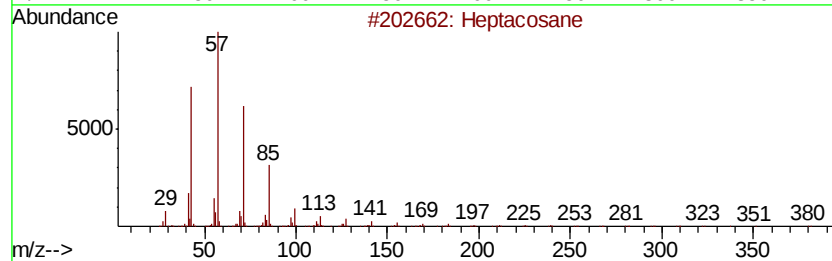
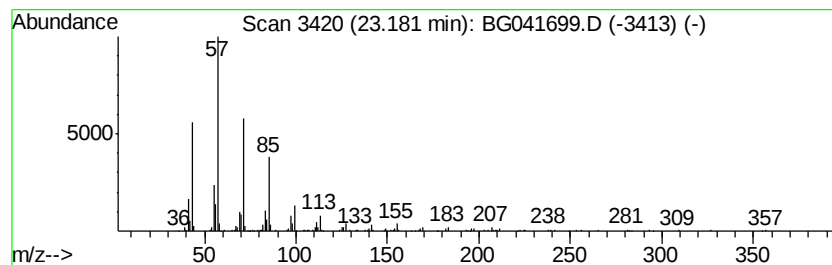
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Heptacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.18	3.35 ng	329582	Chrysene-d12	21.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	91
2		Eicosane	282	C20H42	000112-95-8	91
3		Heptadecane	240	C17H36	000629-78-7	91
4		Heneicosane	296	C21H44	000629-94-7	90
5		Octacosane	394	C28H58	000630-02-4	90



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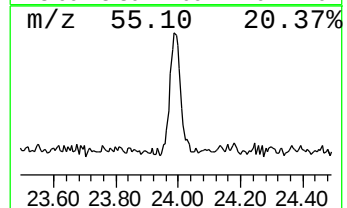
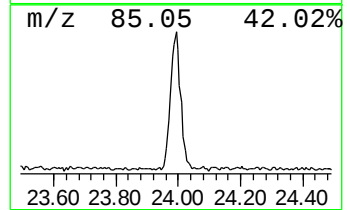
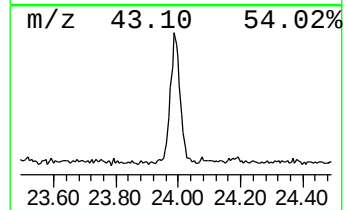
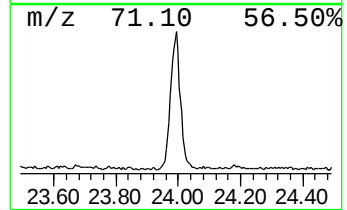
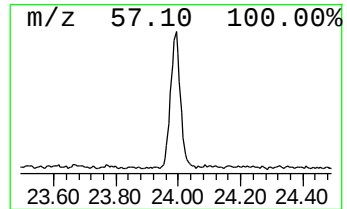
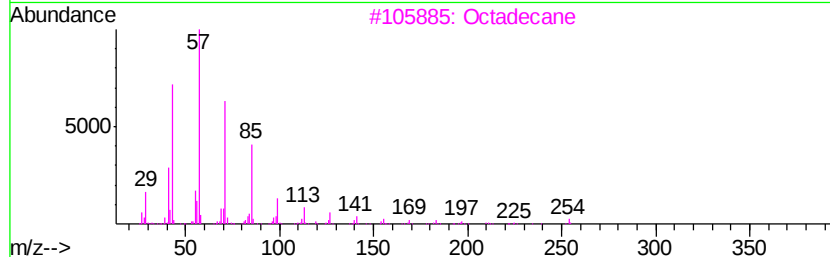
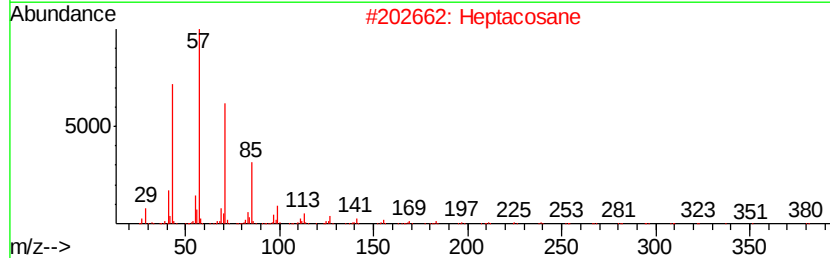
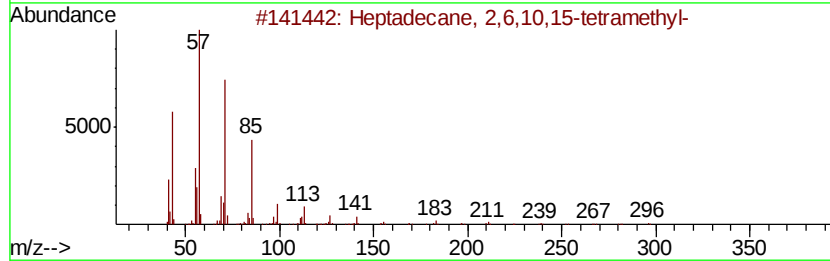
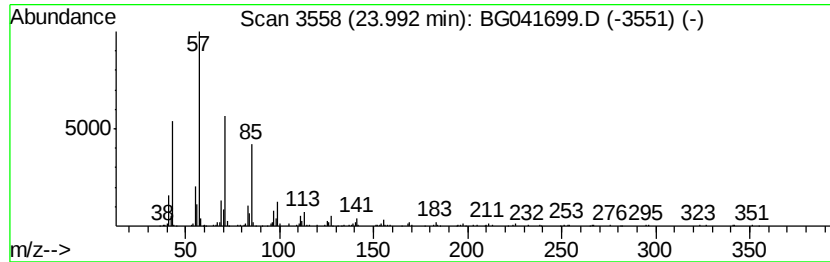
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heptadecane, 2,6,10,15-tetr... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.99	4.12 ng	413433	Perylene-d12	24.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	92
2		Heptacosane	380	C27H56	000593-49-7	91
3		Octadecane	254	C18H38	000593-45-3	91
4		Docosane	310	C22H46	000629-97-0	91
5		Hexadecane	226	C16H34	000544-76-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG071719\
Data File : BG041699.D
Acq On : 17 Jul 2019 22:17
Operator : HP/JU
Sample : K3835-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB-4(7-9)

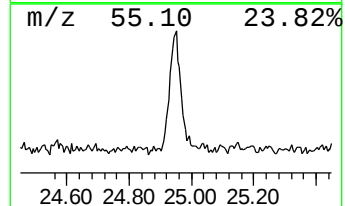
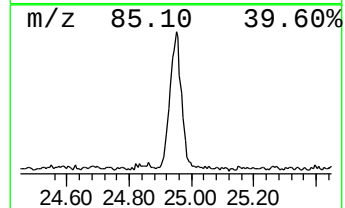
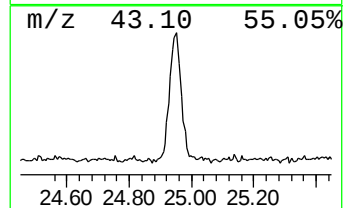
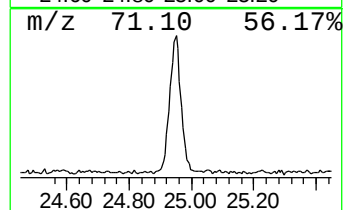
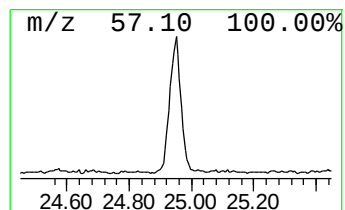
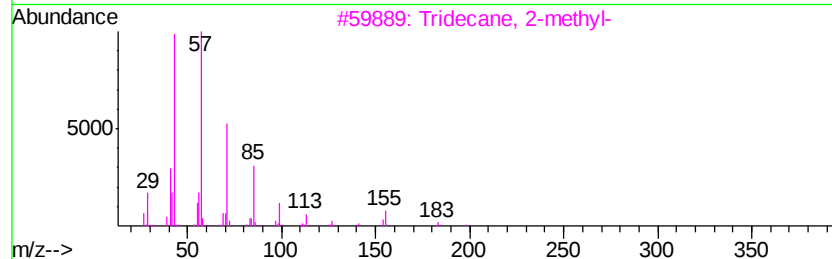
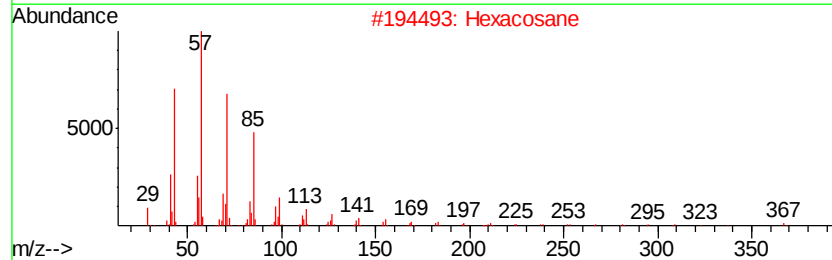
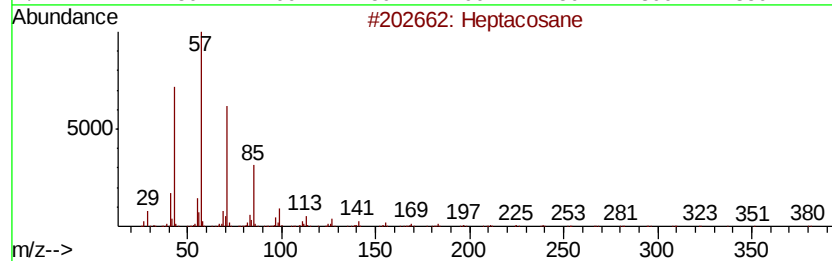
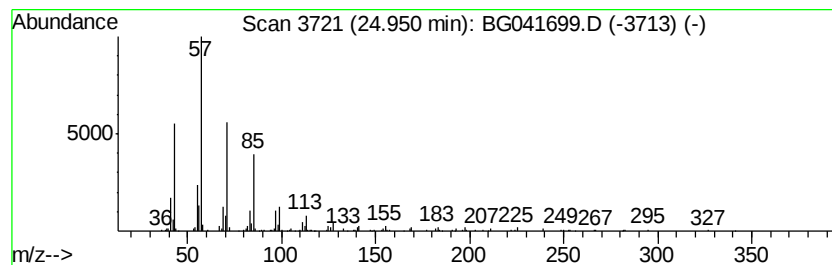
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.95	4.20 ng	420899	Perylene-d12	24.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	91
2		Hexacosane	366	C26H54	000630-01-3	90
3		Tridecane, 2-methyl-	198	C14H30	001560-96-9	90
4		Docosane	310	C22H46	000629-97-0	87
5		Tricosane, 2-methyl-	338	C24H50	001928-30-9	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG071719\
 Data File : BG041699.D
 Acq On : 17 Jul 2019 22:17
 Operator : HP/JU
 Sample : K3835-08
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-4(7-9)

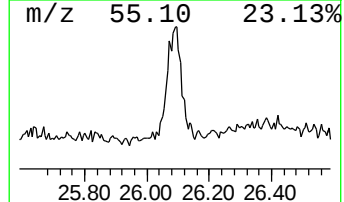
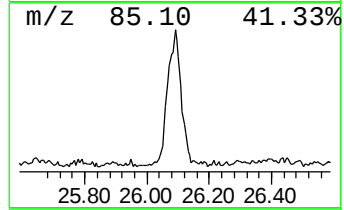
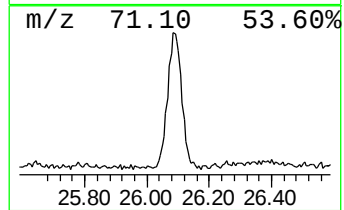
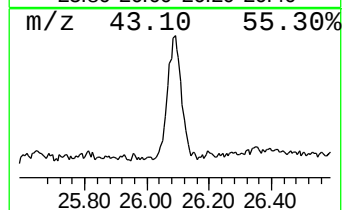
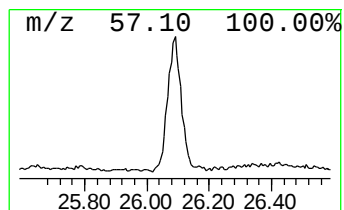
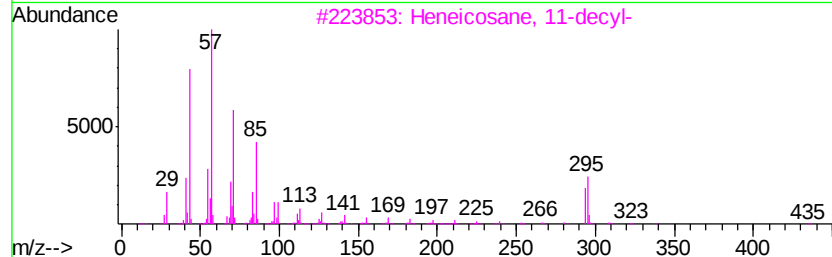
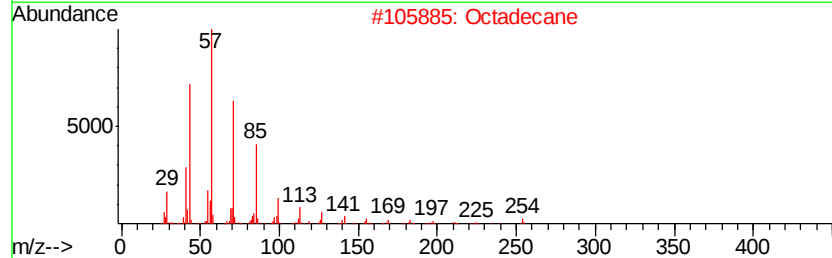
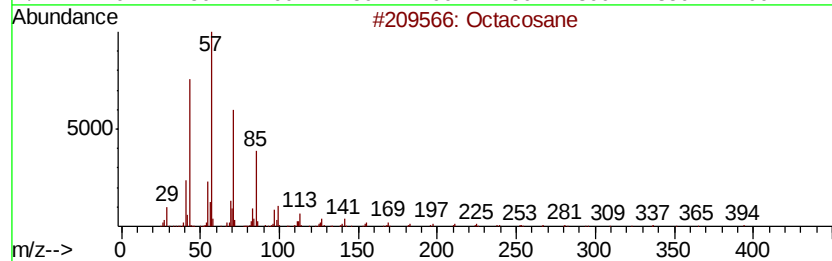
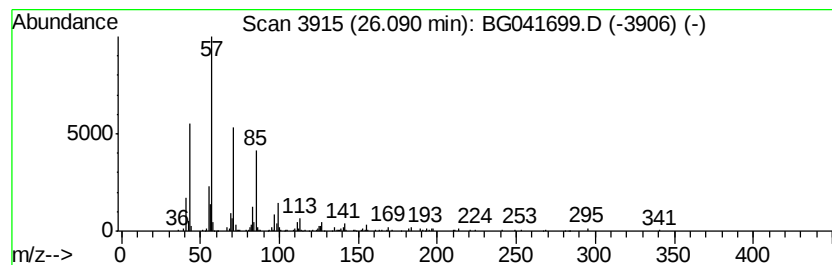
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 9 Heneicosane, 11-decyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.09	3.53 ng	353738	Perylene-d12	24.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Octadecane	254	C18H38	000593-45-3	91
3		Heneicosane, 11-decyl-	437	C31H64	055320-06-4	91
4		Docosane	310	C22H46	000629-97-0	91
5		Eicosane	282	C20H42	000112-95-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG071719\
Data File : BG041699.D
Acq On : 17 Jul 2019 22:17
Operator : HP/JU
Sample : K3835-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB-4(7-9)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG071519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.20	21.3	ng	674751	1	8.05	634597	20.0
unknown7.58	7.58	101.1	ng	3207990	1	8.05	634597	20.0
Tetracosyl heptaf...	21.32	5.7	ng	562567	5	21.65	1969470	20.0
Octacosane	22.48	2.1	ng	203320	5	21.65	1969470	20.0
Heptacosane	23.18	3.4	ng	329582	5	21.65	1969470	20.0
Heptadecane, 2,6,...	23.99	4.1	ng	413433	6	24.84	2005120	20.0
Hexacosane	24.95	4.2	ng	420899	6	24.84	2005120	20.0
Heneicosane, 11-d...	26.09	3.5	ng	353738	6	24.84	2005120	20.0