

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010320\
 Data File : BG043897.D
 Acq On : 3 Jan 2020 21:13
 Operator : JU
 Sample : K6464-06
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 GW-2

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_G\METHODS\8270-BG123019.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	3.143	5	9	26	rVB	1200938	1841728	20.19%	3.680%
2	5.071	331	337	347	rBV	168956	270028	2.96%	0.540%
3	5.535	410	416	426	rBV	974122	1596307	17.50%	3.190%
4	7.121	679	686	699	rBV	668177	1240946	13.60%	2.480%
5	7.497	742	750	758	rBV	1956350	3389010	37.15%	6.772%
6	7.961	822	829	837	rBV	446625	753732	8.26%	1.506%
7	8.284	876	884	894	rBV	1734950	3024485	33.16%	6.044%
8	9.119	1017	1026	1034	rBV	1441665	2662941	29.19%	5.321%
9	10.764	1297	1306	1314	rBV	645874	1147700	12.58%	2.293%
10	13.220	1716	1724	1731	rBV	3969519	6192296	67.88%	12.374%
11	14.042	1853	1864	1869	rBV	106956	162077	1.78%	0.324%
12	14.595	1949	1958	1966	rBV2	1074348	1722238	18.88%	3.441%
13	16.081	2204	2211	2228	rBV	3329695	5167719	56.65%	10.326%
14	17.333	2418	2424	2431	rBV2	1363881	2097085	22.99%	4.190%
15	18.237	2574	2578	2586	rBV4	54229	92198	1.01%	0.184%
16	19.953	2863	2870	2880	rBV	6475605	9122116	100.00%	18.228%
17	21.257	3087	3092	3101	rBV2	634105	1023365	11.22%	2.045%
18	21.586	3141	3148	3159	rVB	1381794	2239331	24.55%	4.475%
19	21.810	3181	3186	3193	rVB	189323	270902	2.97%	0.541%
20	22.409	3282	3288	3300	rVB	281300	499951	5.48%	0.999%
21	23.085	3397	3403	3414	rBV	355702	655736	7.19%	1.310%
22	23.878	3530	3538	3551	rBV	361623	804307	8.82%	1.607%
23	24.712	3670	3680	3689	rBV	821779	2244978	24.61%	4.486%
24	24.806	3689	3696	3705	rVB2	322484	766744	8.41%	1.532%
25	25.911	3875	3884	3894	rBV3	221737	635759	6.97%	1.270%
26	26.011	3899	3901	3913	rVB10	40075	96566	1.06%	0.193%
27	27.233	4101	4109	4121	rVB5	95677	323677	3.55%	0.647%

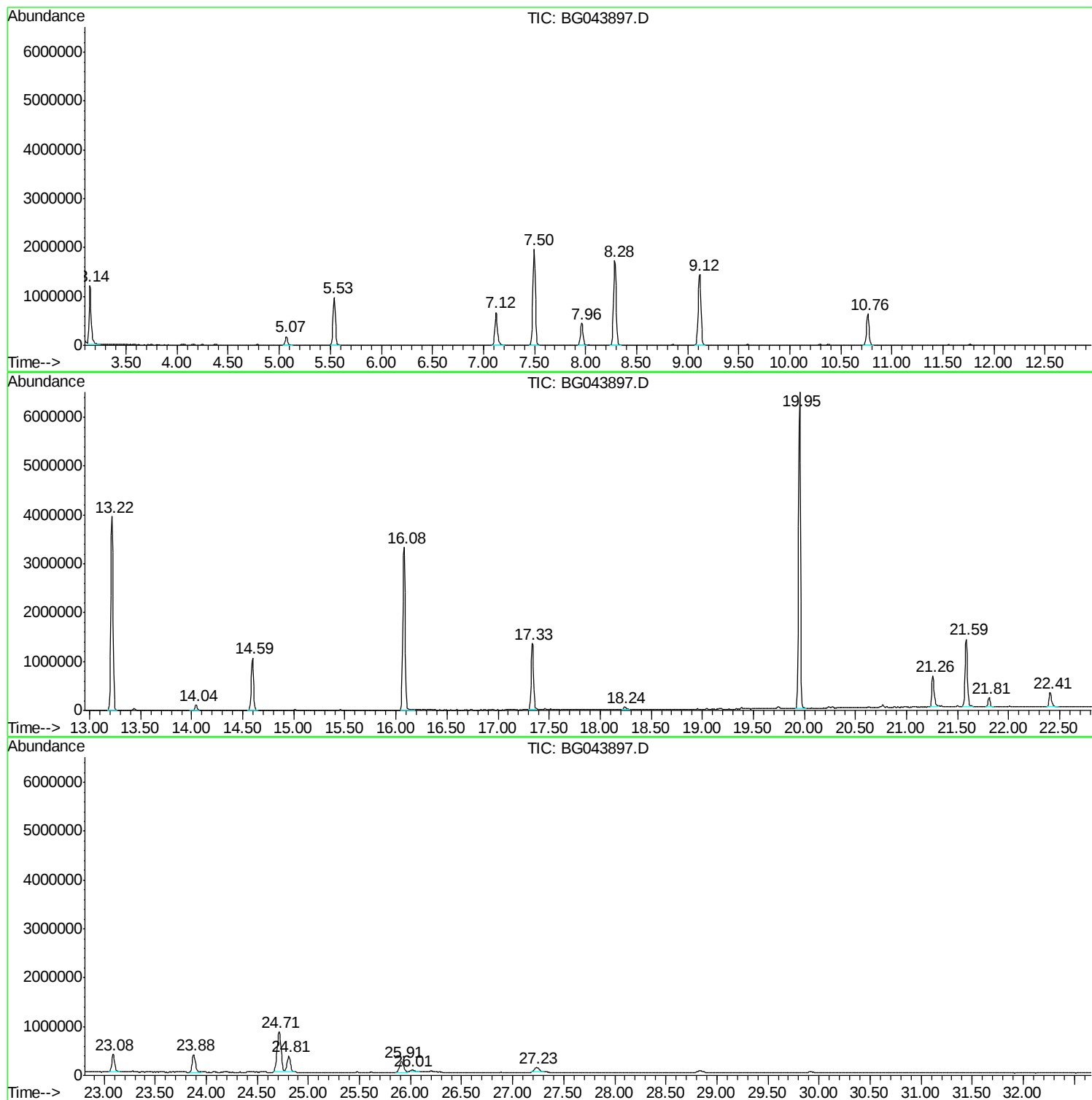
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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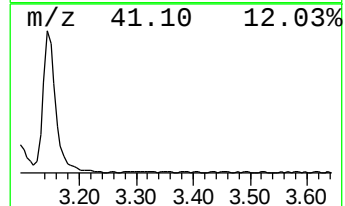
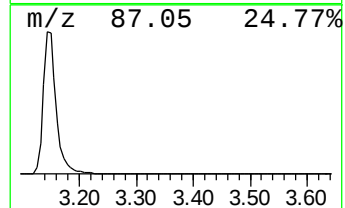
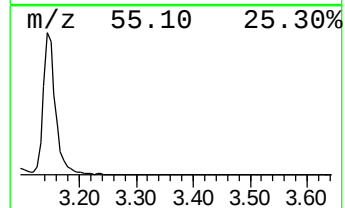
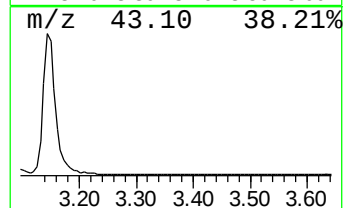
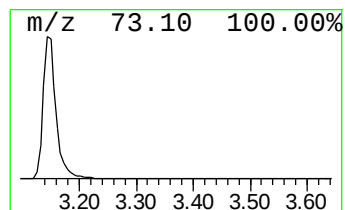
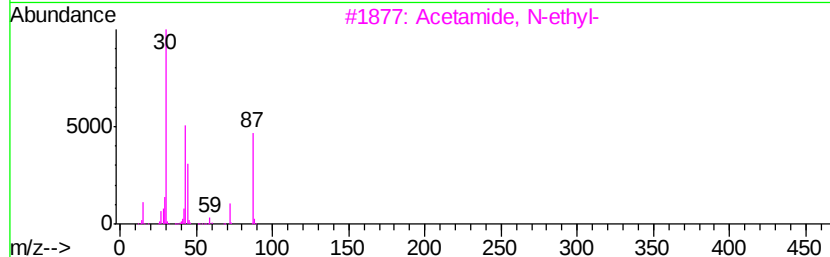
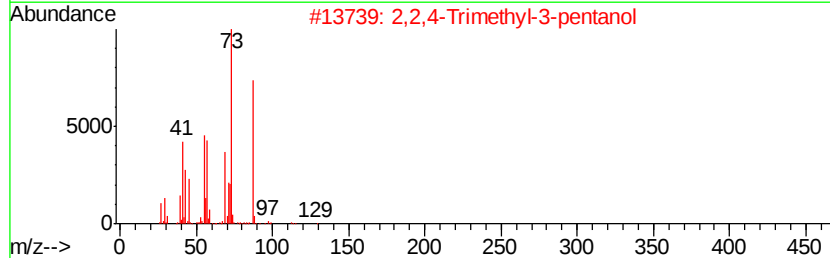
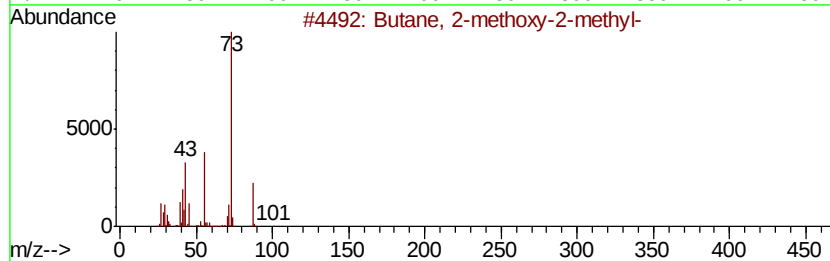
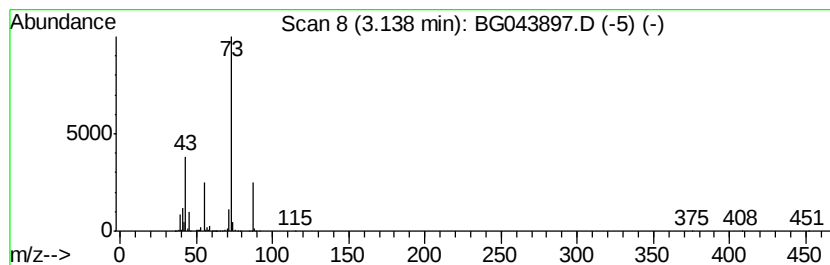
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
3.14	48.87 ng	1841730	1,4-Dichlorobenzene-d4	7.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	38
3		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	35
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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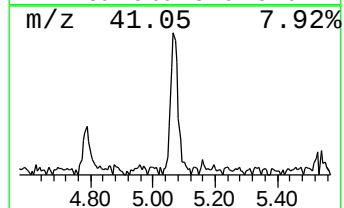
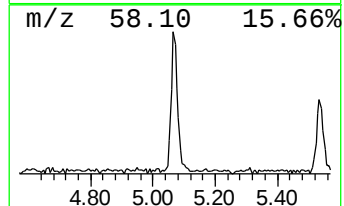
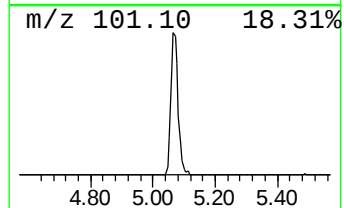
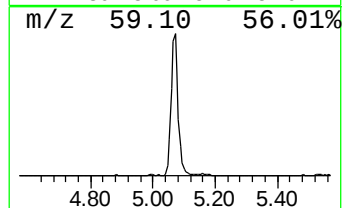
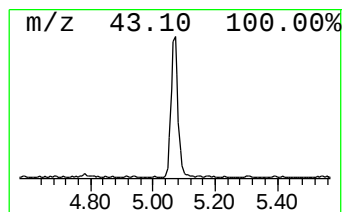
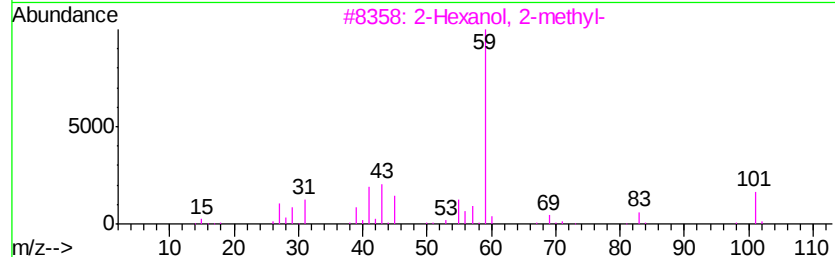
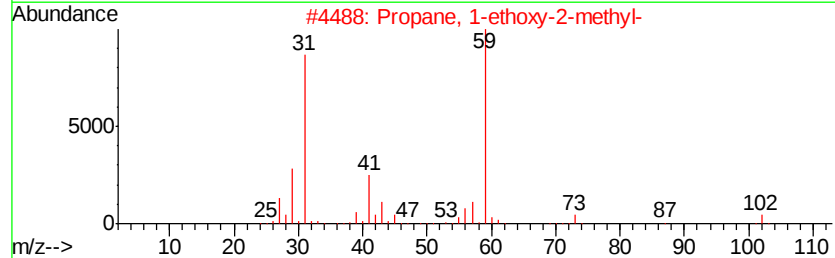
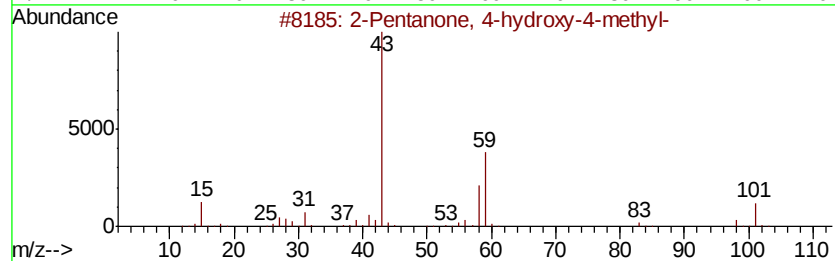
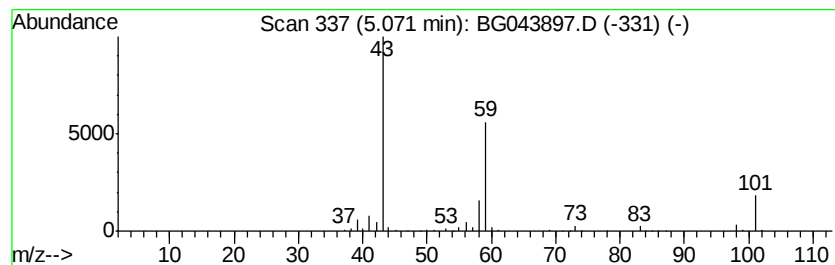
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	7.17 ng	270028	1,4-Dichlorobenzene-d4	7.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	32
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
4			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16



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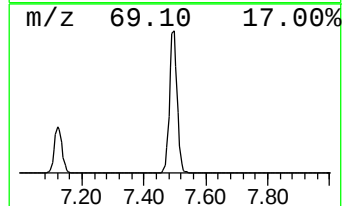
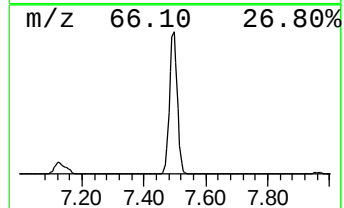
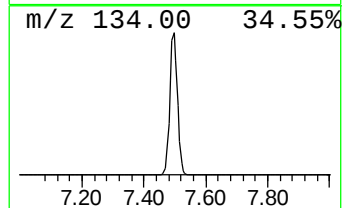
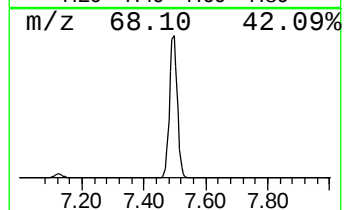
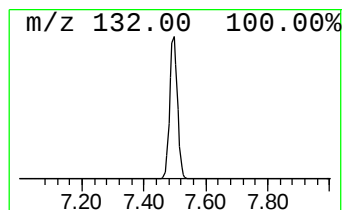
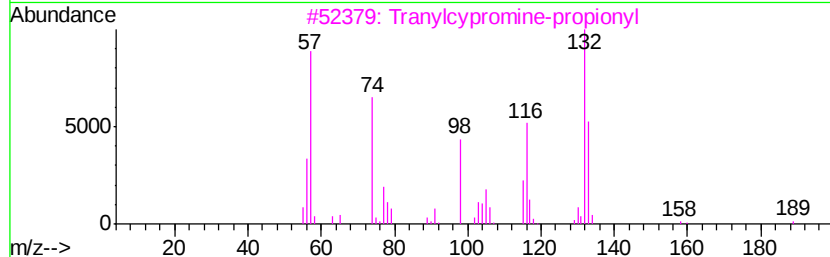
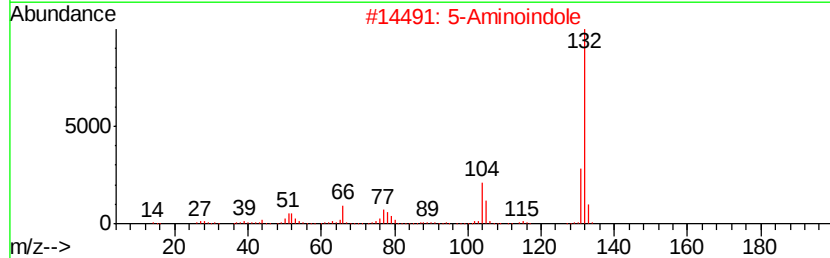
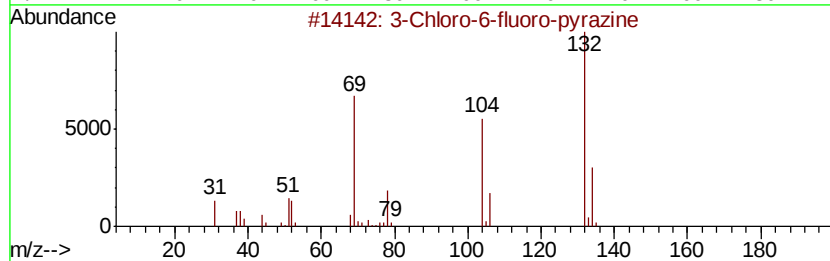
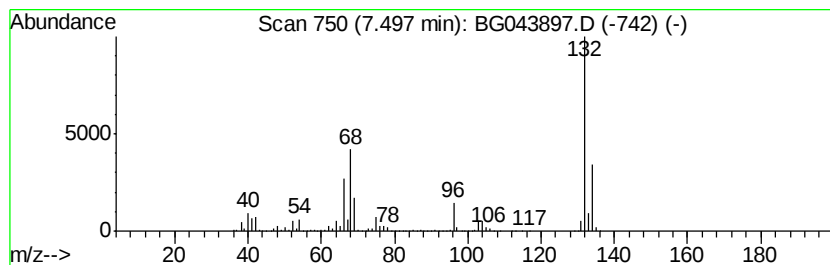
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Peak Number 3 unknown7.50 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.50	89.93 ng	3389010	1,4-Dichlorobenzene-d4	7.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	22
3			Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
5			Xenon	132	Xe	007440-63-3	9



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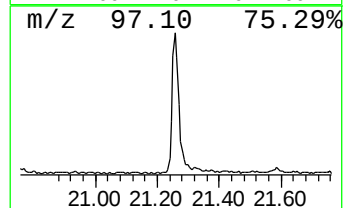
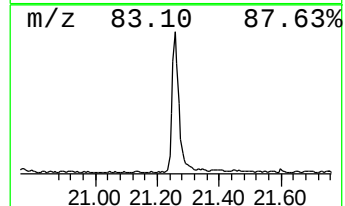
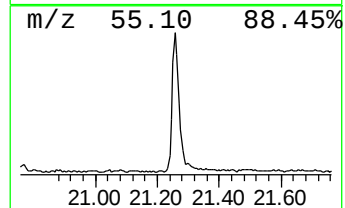
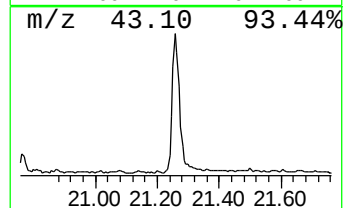
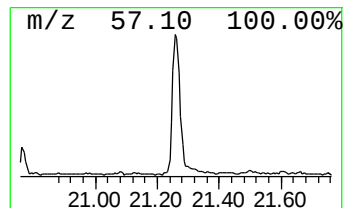
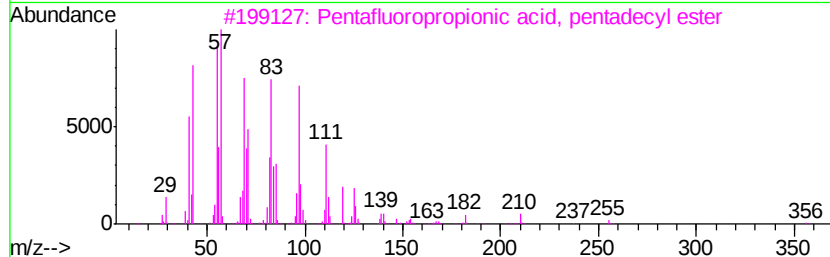
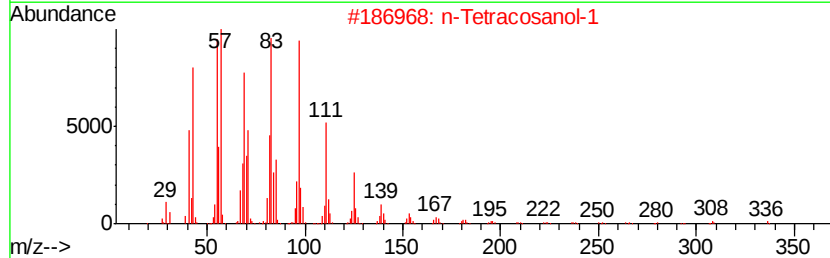
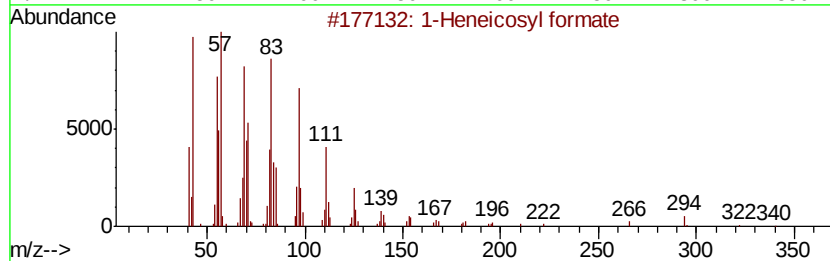
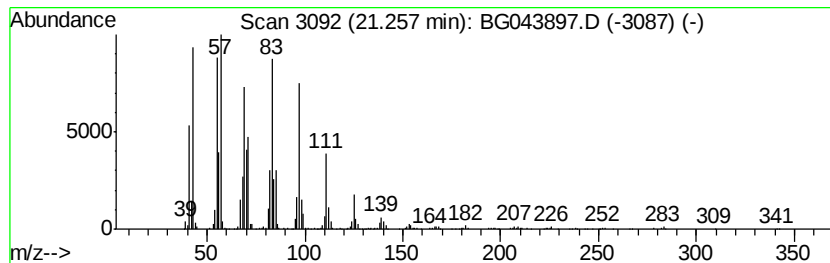
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Peak Number 5 1-Heneicosyl formate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.26	9.14 ng	1023370	Chrysene-d12	21.59

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
2		n-Tetracosanol-1	354	C24H50O	000506-51-4	94
3		Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	93
4		Cyclotridecane	182	C13H26	000295-02-3	92
5		Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	91



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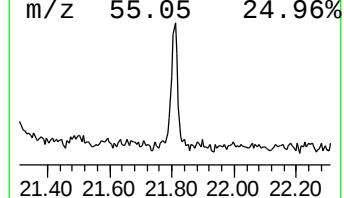
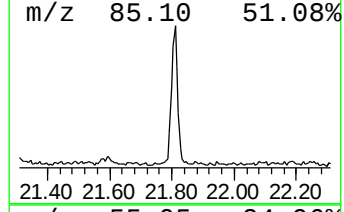
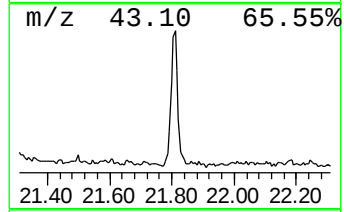
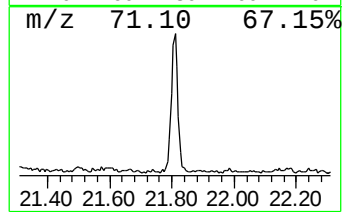
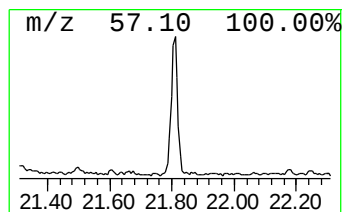
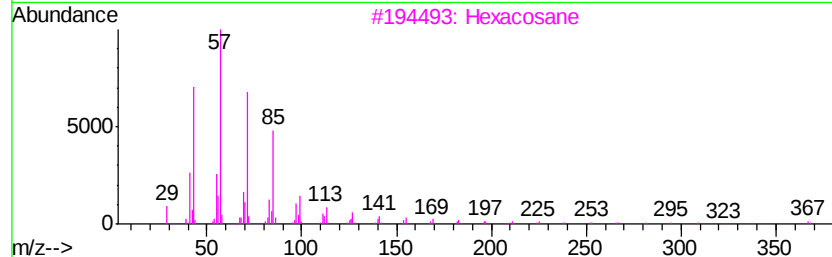
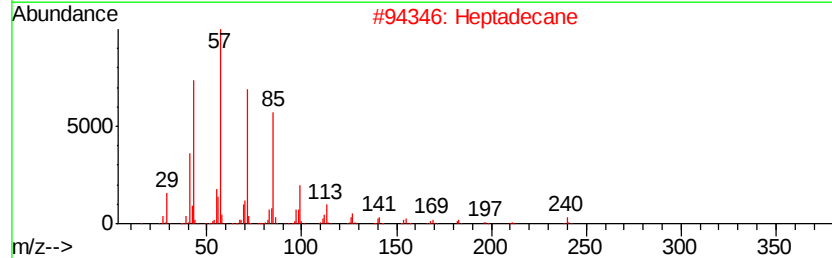
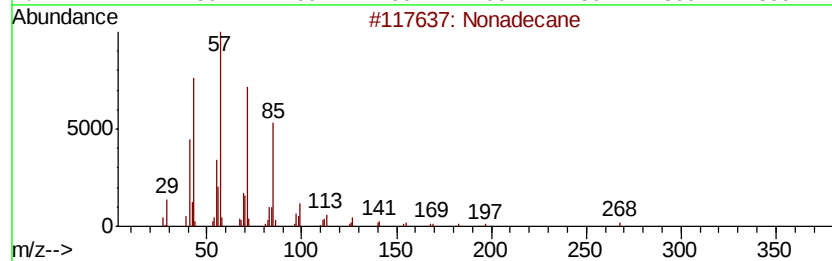
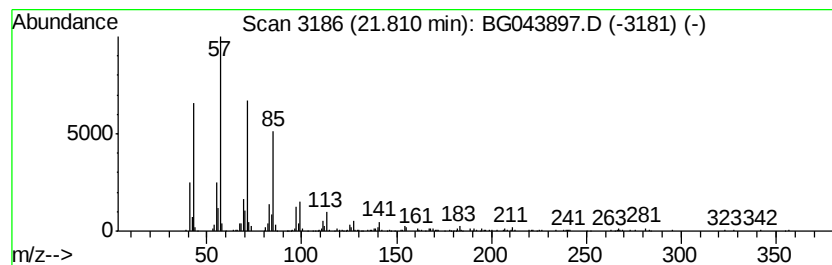
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Peak Number 6 Nonadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.81	2.42 ng	270902	Chrysene-d12	21.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	95
2		Heptadecane	240	C17H36	000629-78-7	93
3		Hexacosane	366	C26H54	000630-01-3	91
4		Octacosane	394	C28H58	000630-02-4	91
5		Heneicosane	296	C21H44	000629-94-7	91



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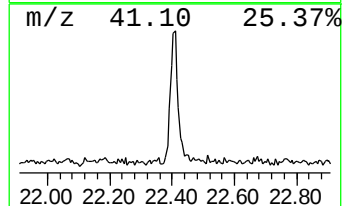
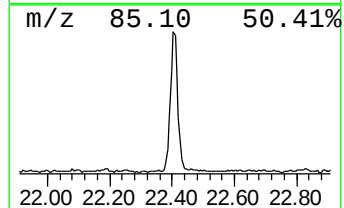
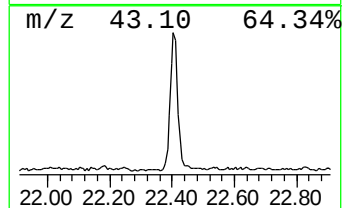
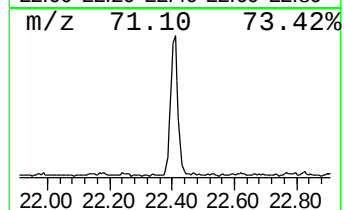
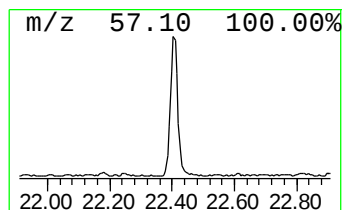
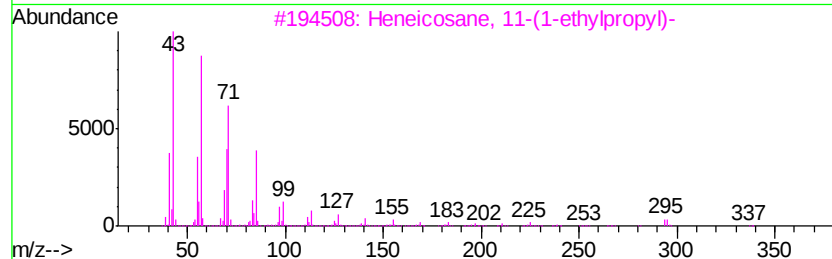
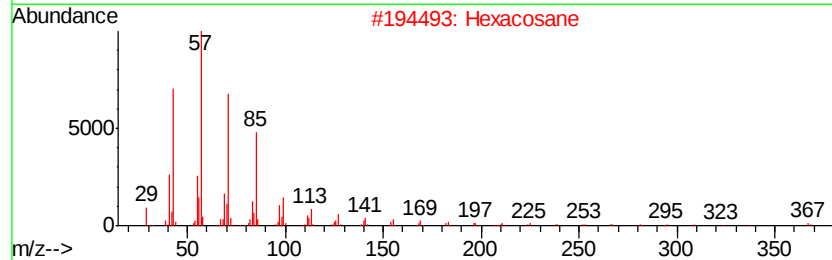
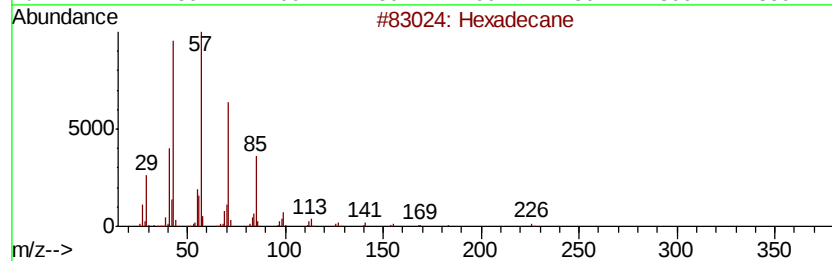
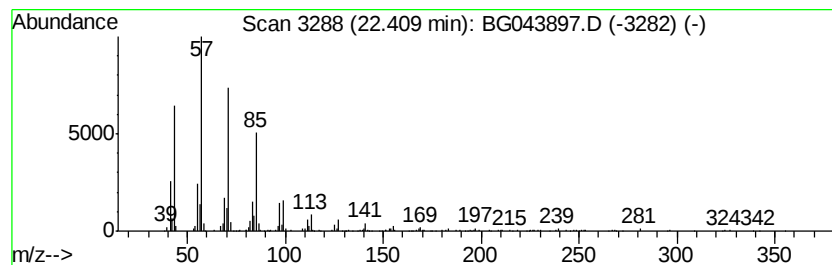
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.41	4.47 ng	499951	Chrysene-d12	21.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	93
2		Hexacosane	366	C26H54	000630-01-3	91
3		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
4		Triacontane	422	C30H62	000638-68-6	91
5		Tricosane	324	C23H48	000638-67-5	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010320\
 Data File : BG043897.D
 Acq On : 3 Jan 2020 21:13
 Operator : JU
 Sample : K6464-06
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 GW-2

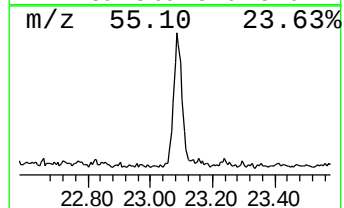
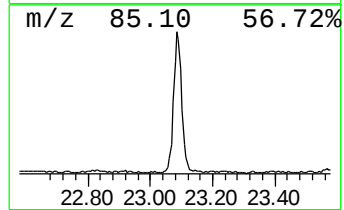
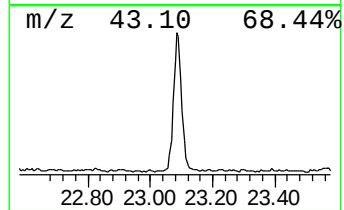
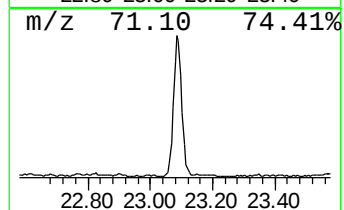
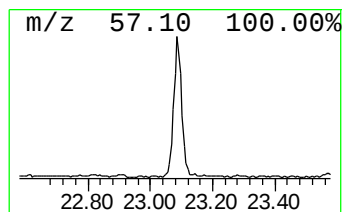
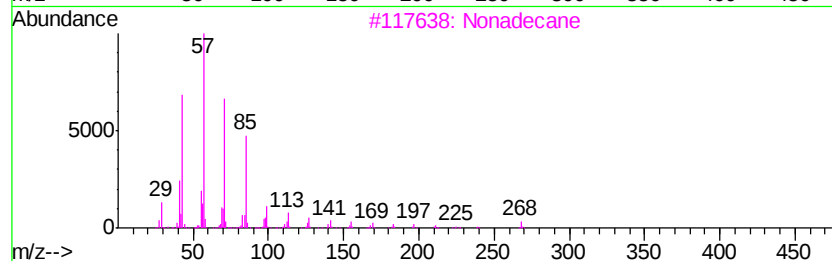
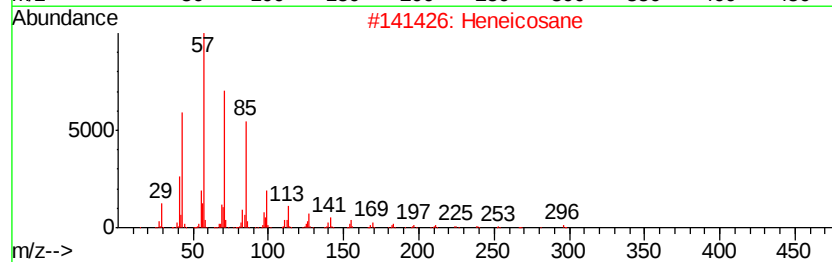
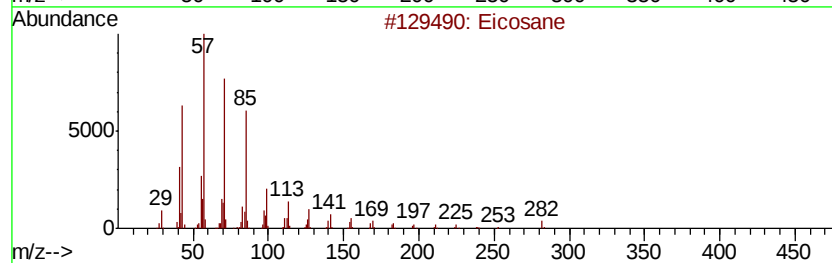
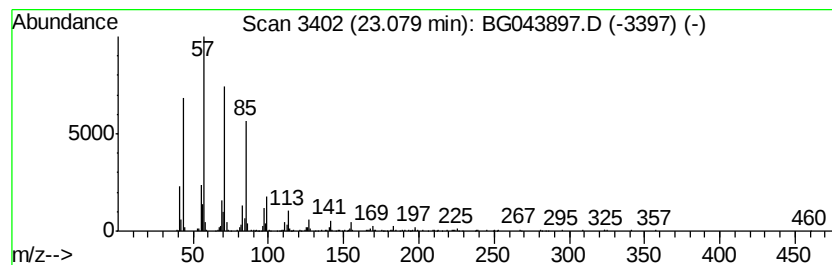
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 8 Eicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.08	5.86 ng	655736	Chrysene-d12	21.59

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	98
2	Heneicosane		296	C21H44	000629-94-7	90
3	Nonadecane		268	C19H40	000629-92-5	86
4	Tetratetracontane		619	C44H90	007098-22-8	83
5	Hexacosane		366	C26H54	000630-01-3	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010320\
Data File : BG043897.D
Acq On : 3 Jan 2020 21:13
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Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GW-2

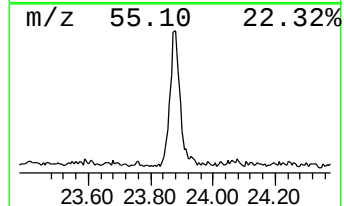
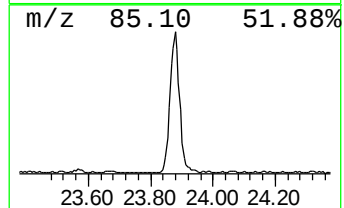
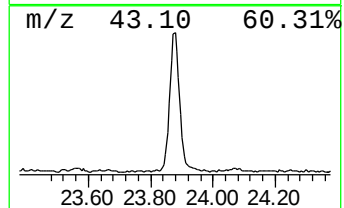
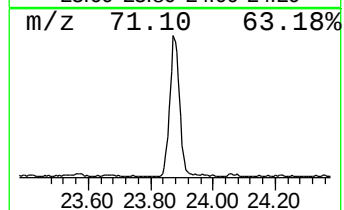
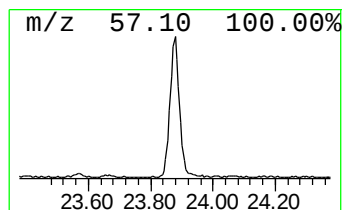
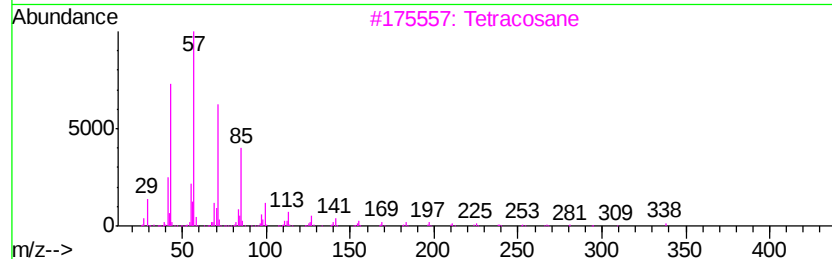
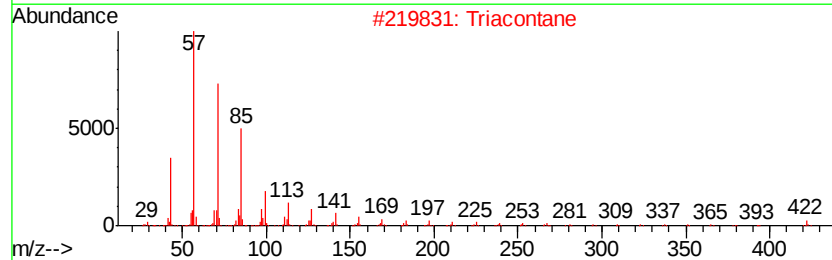
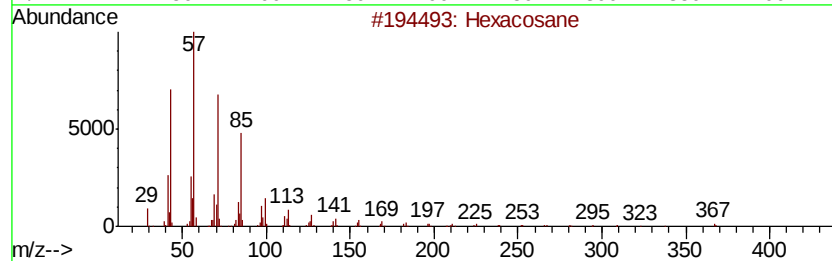
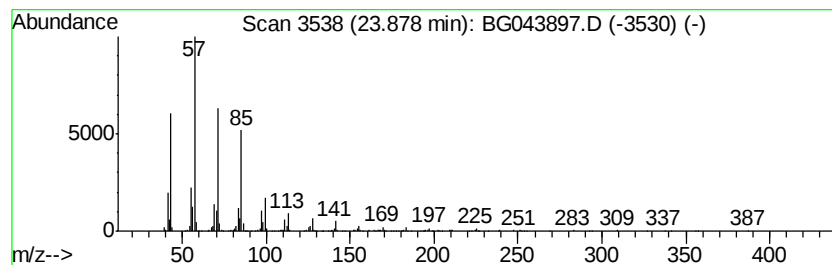
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.88	7.17 ng	804307	Perylene-d12	24.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	91
2		Triacontane	422	C30H62	000638-68-6	91
3		Tetracosane	338	C24H50	000646-31-1	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Eicosane	282	C20H42	000112-95-8	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010320\
 Data File : BG043897.D
 Acq On : 3 Jan 2020 21:13
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 Sample : K6464-06
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Instrument :
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 ClientSampleId :
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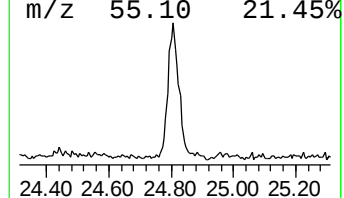
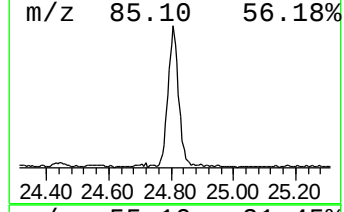
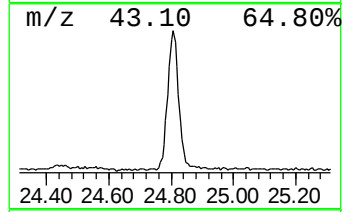
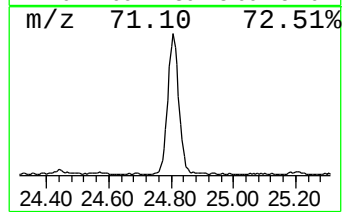
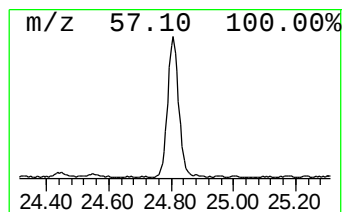
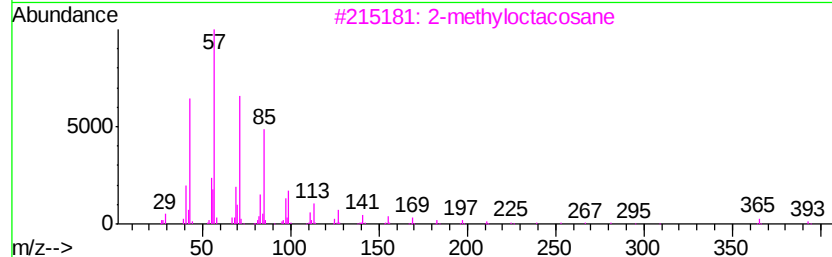
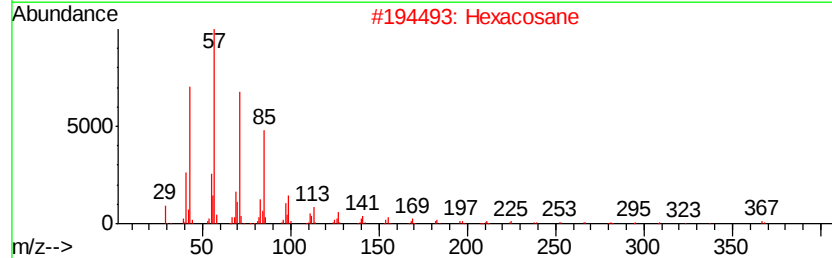
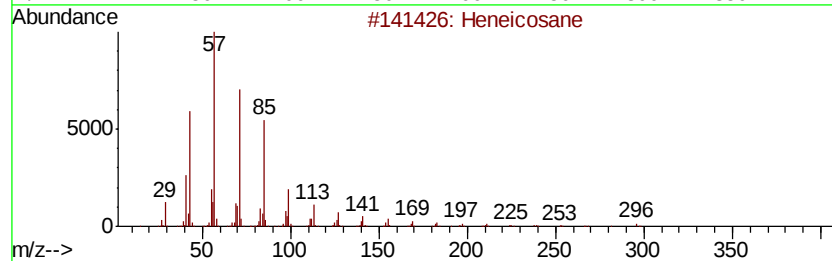
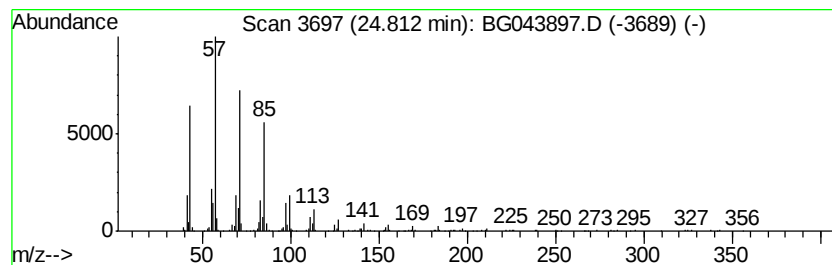
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 10 Heneicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.81	6.83 ng	766744	Perylene-d12	24.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	91
2			Hexacosane	366	C26H54	000630-01-3	91
3			2-methyloctacosane	408	C29H60	1000376-72-8	91
4			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5			Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010320\
 Data File : BG043897.D
 Acq On : 3 Jan 2020 21:13
 Operator : JU
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 Misc :
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 ClientSampleId :
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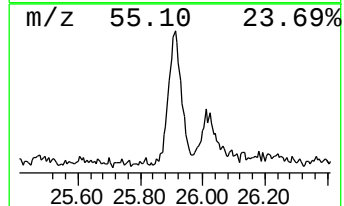
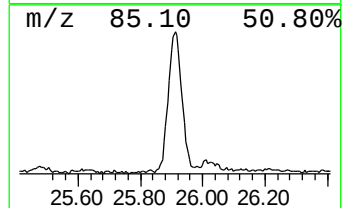
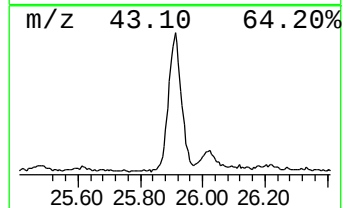
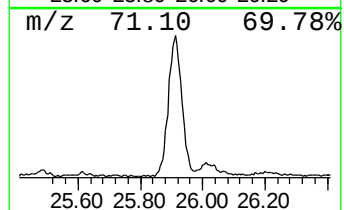
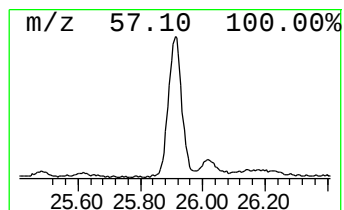
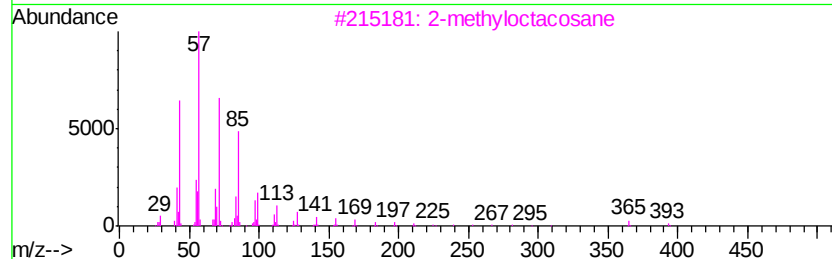
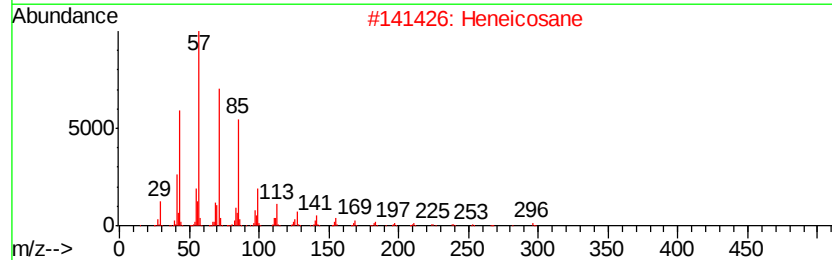
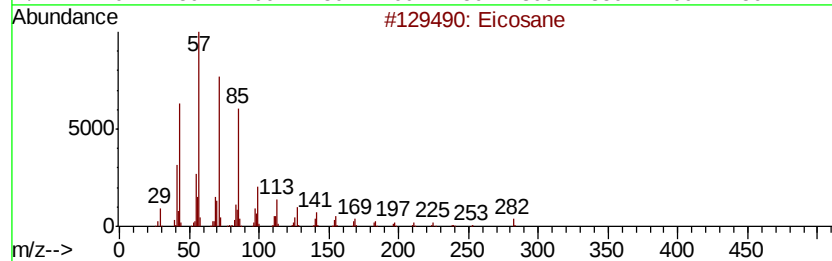
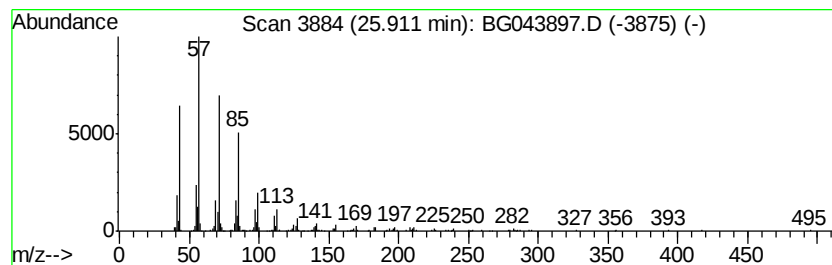
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 11 2-methyloctacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.91	5.66 ng	635759	Perylene-d12	24.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	95
2			Heneicosane	296	C21H44	000629-94-7	90
3			2-methyloctacosane	408	C29H60	1000376-72-8	87
4			Hexacosane	366	C26H54	000630-01-3	87
5			Heneicosane, 11-decyl-	437	C31H64	055320-06-4	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010320\
Data File : BG043897.D
Acq On : 3 Jan 2020 21:13
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Misc :
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Instrument :
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ClientSampleId :
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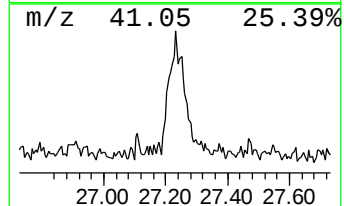
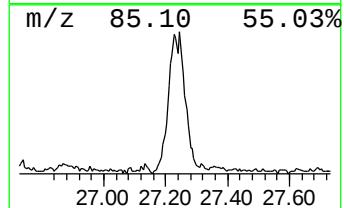
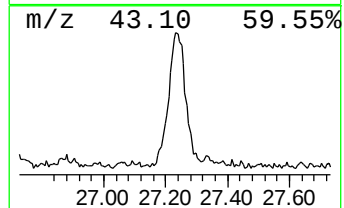
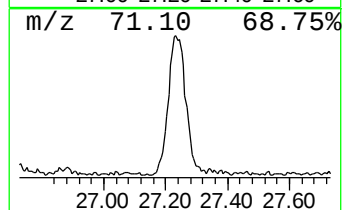
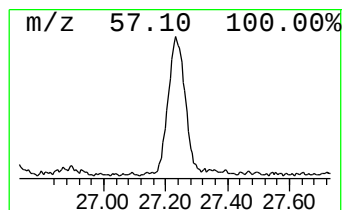
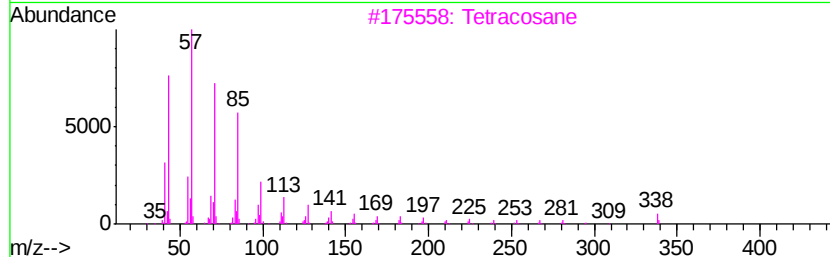
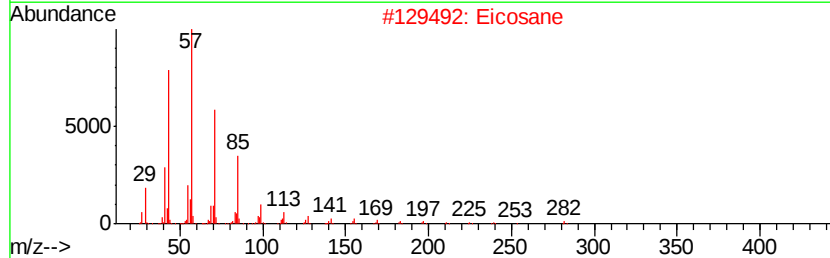
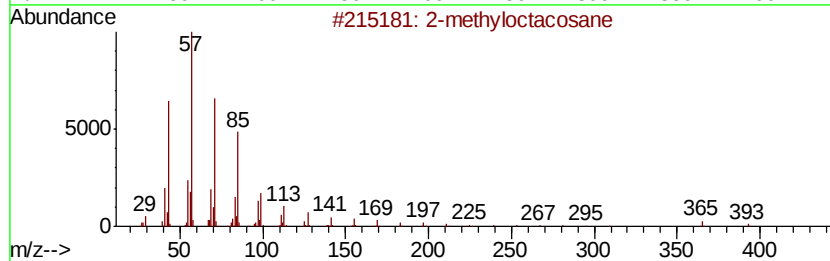
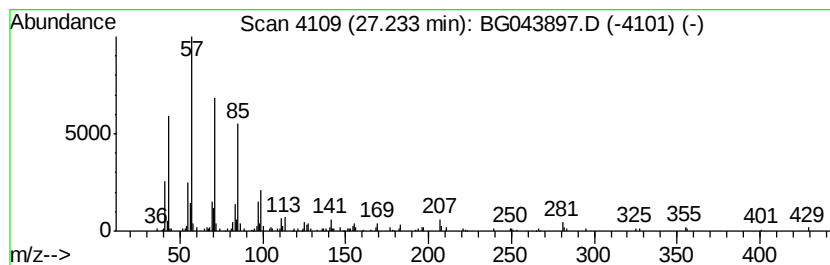
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tetracosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.23	2.88 ng	323677	Perylene-d12	24.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-methyloctacosane	408	C29H60	1000376-72-8	86
2			Eicosane	282	C20H42	000112-95-8	83
3			Tetracosane	338	C24H50	000646-31-1	80
4			Octacosane	394	C28H58	000630-02-4	74
5			Hexacosane	366	C26H54	000630-01-3	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG010320\
Data File : BG043897.D
Acq On : 3 Jan 2020 21:13
Operator : JU
Sample : K6464-06
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GW-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG123019.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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2-Pentanone, 4-hy...	5.07	7.2	ng	270028	1	7.96	753732	20.0
unknown7.50	7.50	89.9	ng	3389010	1	7.96	753732	20.0
1-Heneicosyl formate	21.26	9.1	ng	1023370	5	21.59	2239330	20.0
Nonadecane	21.81	2.4	ng	270902	5	21.59	2239330	20.0
Hexadecane	22.41	4.5	ng	499951	5	21.59	2239330	20.0
Eicosane	23.08	5.9	ng	655736	5	21.59	2239330	20.0
Hexacosane	23.88	7.2	ng	804307	6	24.71	2244980	20.0
Heneicosane	24.81	6.8	ng	766744	6	24.71	2244980	20.0
2-methyloctacosane	25.91	5.7	ng	635759	6	24.71	2244980	20.0
Tetracosane	27.23	2.9	ng	323677	6	24.71	2244980	20.0