

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082619\
 Data File : BG042502.D
 Acq On : 26 Aug 2019 17:10
 Operator : HP/JU
 Sample : K4491-06
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WP-1

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-BG082219.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	3.118	3	5	34	rVB	1533969	2861309	99.72%	14.798%
2	5.556	414	420	438	rBV	204546	366422	12.77%	1.895%
3	7.166	684	694	709	rBV	119446	246650	8.60%	1.276%
4	7.530	749	756	769	rBV	411376	745261	25.97%	3.854%
5	8.000	829	836	844	rBV	114519	193688	6.75%	1.002%
6	8.323	884	891	901	rBV	447546	792261	27.61%	4.097%
7	9.164	1024	1034	1047	rBV	278039	538381	18.76%	2.784%
8	10.820	1307	1316	1322	rBV	135169	249416	8.69%	1.290%
9	13.276	1727	1734	1747	rBV	945691	1490304	51.94%	7.708%
10	14.105	1869	1875	1879	rBV	24434	39180	1.37%	0.203%
11	14.657	1960	1969	1976	rBV2	223540	364604	12.71%	1.886%
12	16.149	2215	2223	2233	rBV	712789	1175338	40.96%	6.079%
13	16.390	2258	2264	2269	rBV3	17888	29707	1.04%	0.154%
14	17.413	2432	2438	2442	rBV2	294046	453310	15.80%	2.344%
15	17.454	2442	2445	2451	rVB	55231	78459	2.73%	0.406%
16	18.312	2576	2591	2607	rBV	1053361	2122076	73.96%	10.975%
17	18.846	2678	2682	2685	rVB3	23754	31506	1.10%	0.163%
18	19.463	2778	2787	2793	rBV	66187	150860	5.26%	0.780%
19	19.581	2800	2807	2817	rBV	616027	987355	34.41%	5.106%
20	19.828	2846	2849	2854	rVB	41719	56959	1.99%	0.295%
21	20.021	2876	2882	2888	rBV	2145537	2869217	100.00%	14.839%
22	20.327	2931	2934	2938	rVB3	32080	41472	1.45%	0.214%
23	20.820	3015	3018	3023	rVB2	43826	57141	1.99%	0.296%
24	21.332	3101	3105	3116	rVB2	66813	130247	4.54%	0.674%
25	21.690	3159	3166	3172	rBV	448683	784947	27.36%	4.060%
26	21.878	3194	3198	3207	rVB	85469	126493	4.41%	0.654%
27	22.495	3298	3303	3310	rBV	100272	159729	5.57%	0.826%
28	23.194	3416	3422	3430	rVB	120961	226199	7.88%	1.170%
29	23.382	3448	3454	3462	rVB	125191	255677	8.91%	1.322%
30	24.005	3554	3560	3570	rVB2	106730	242794	8.46%	1.256%
31	24.927	3708	3717	3748	rVB3	322500	1264772	44.08%	6.541%
32	26.108	3911	3918	3930	rVB3	71043	203888	7.11%	1.054%

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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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

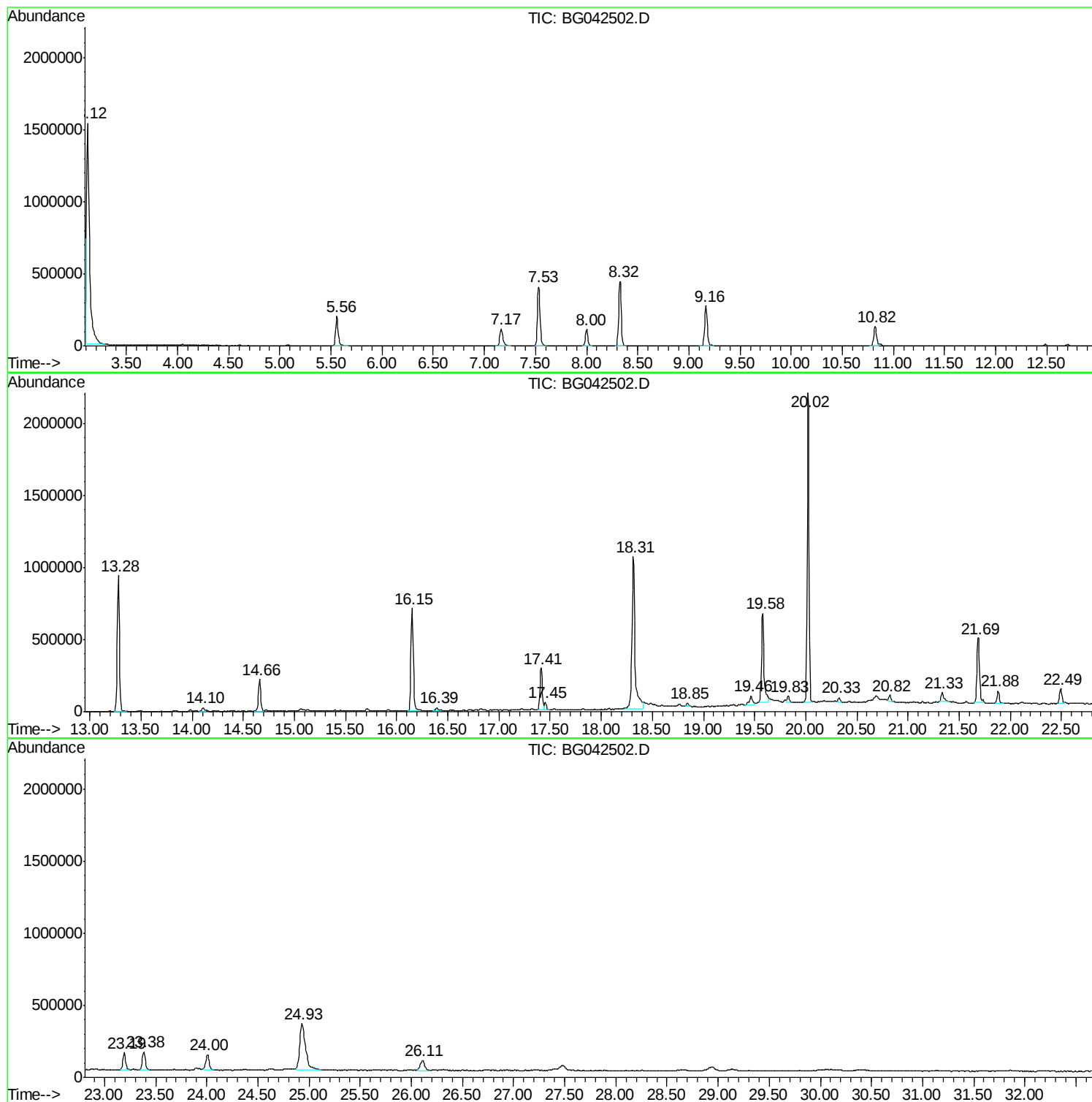
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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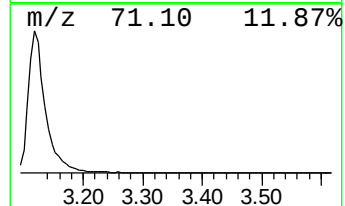
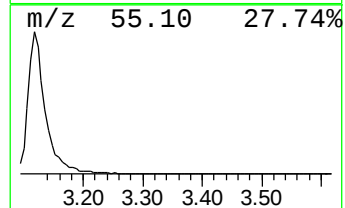
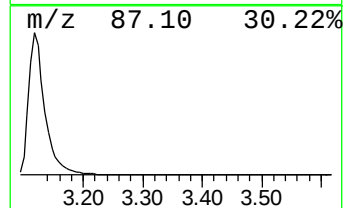
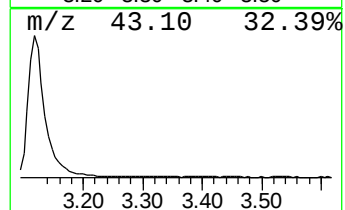
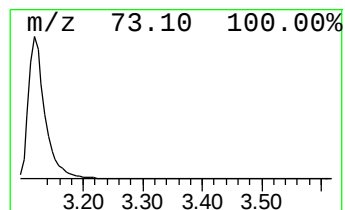
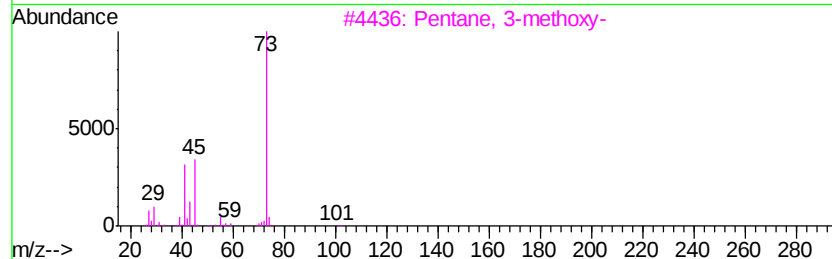
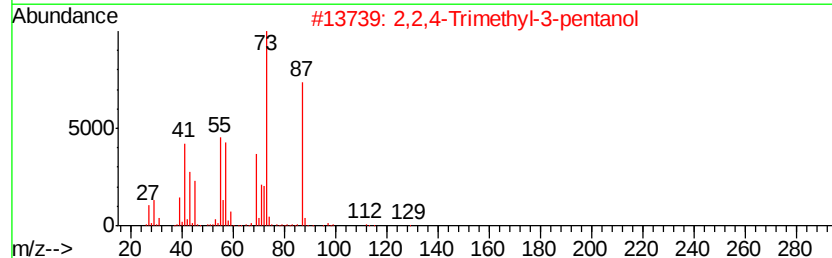
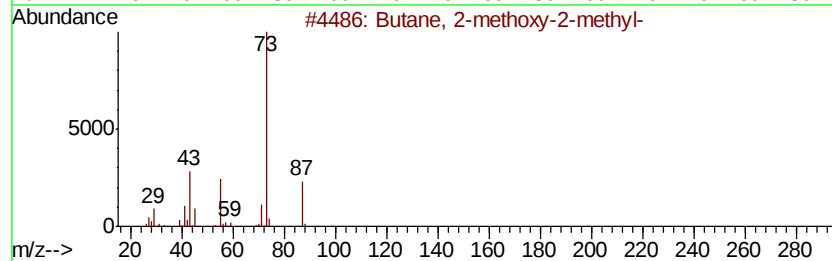
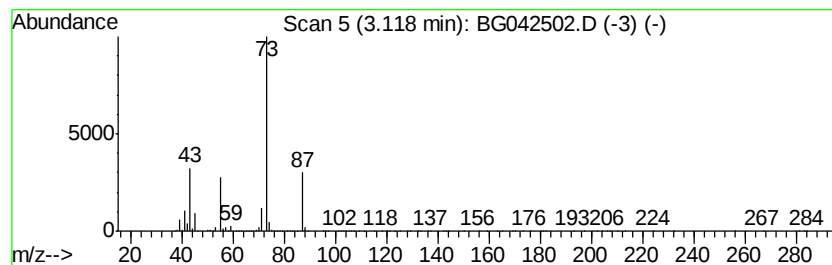
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.12	295.46 ng	2861310	1,4-Dichlorobenzene-d4	8.00

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	40
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
5		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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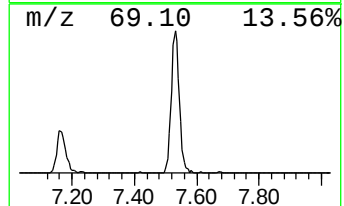
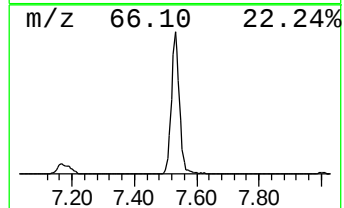
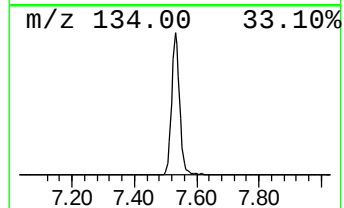
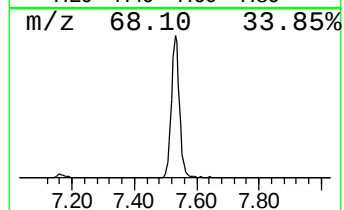
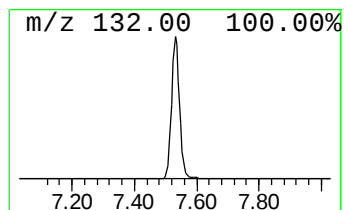
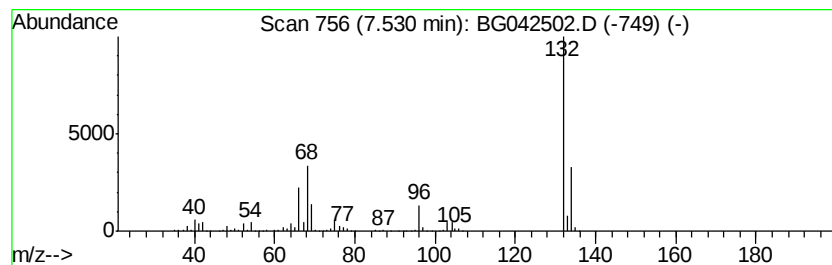
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TIC Library : C:\Database\NIST11.L
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Peak Number 2 unknown7.53 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.53	76.95 ng	745261	1,4-Dichlorobenzene-d4	8.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	12
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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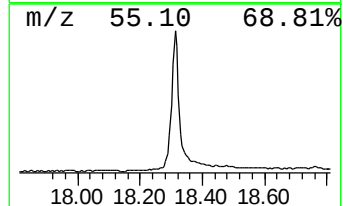
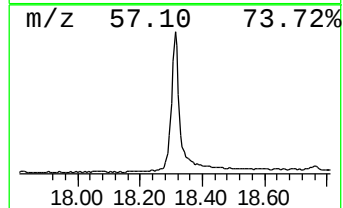
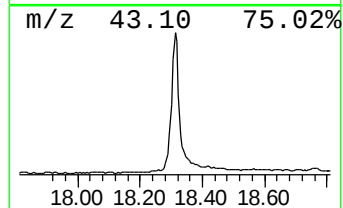
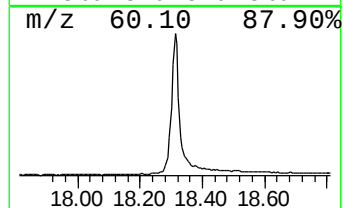
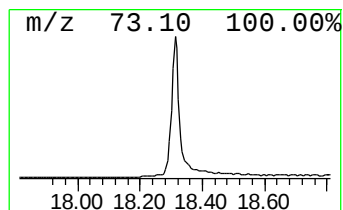
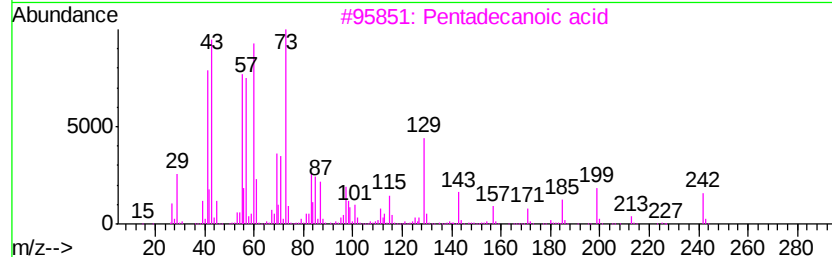
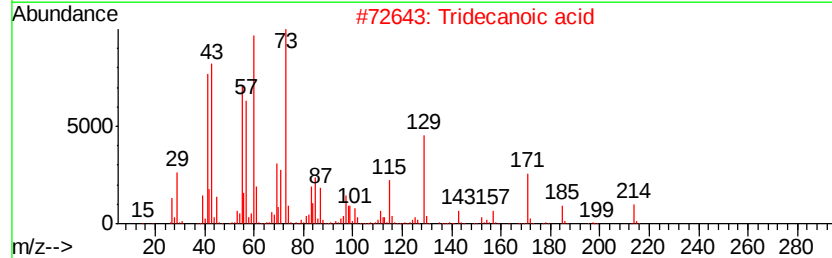
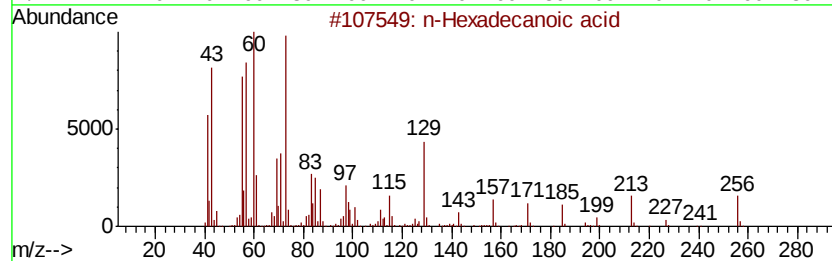
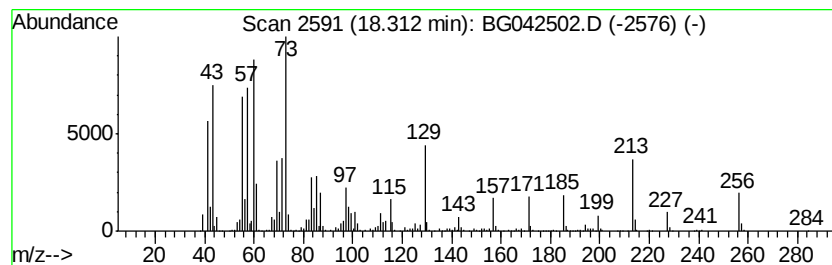
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TIC Library : C:\Database\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 4 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.31	93.63 ng	2122080	Phenanthrene-d10	17.41

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	89
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	81
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	64
5		Undecanoic acid	186	C11H22O2	000112-37-8	59



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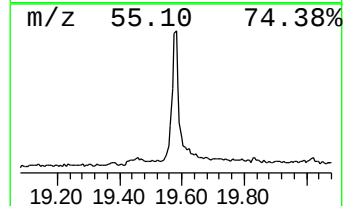
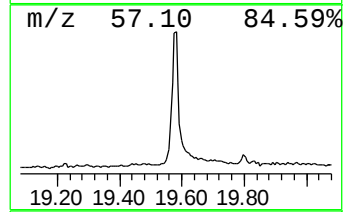
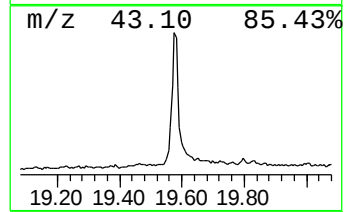
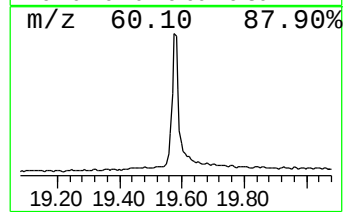
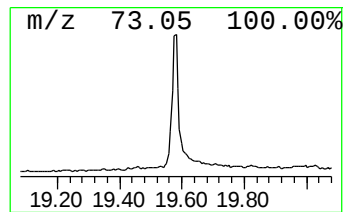
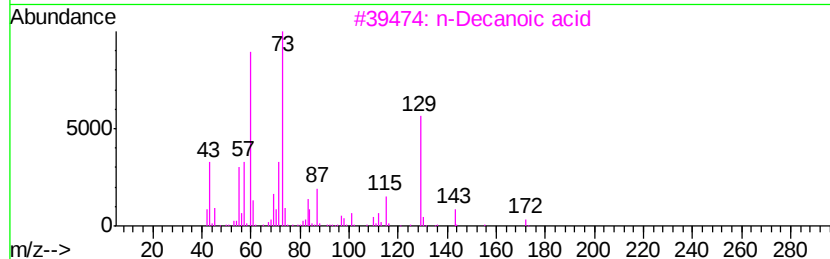
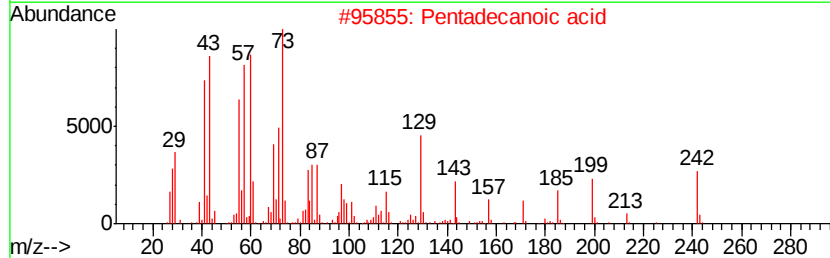
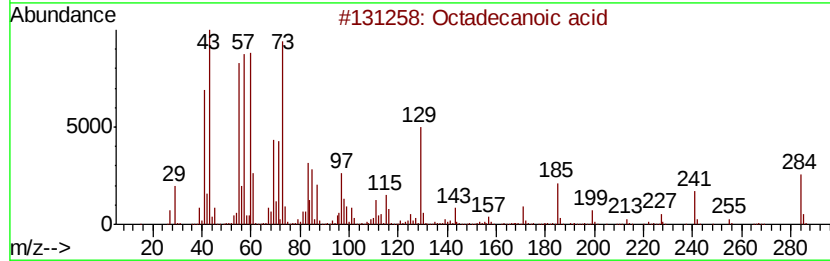
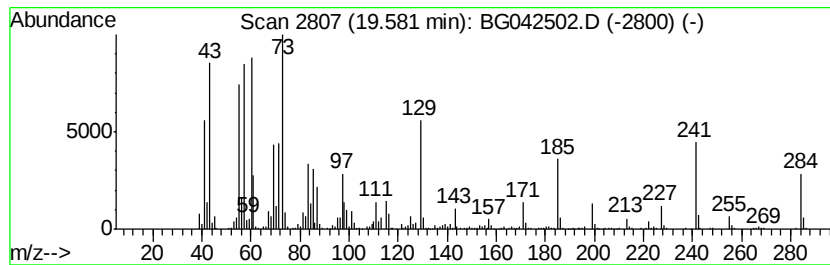
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TIC Library : C:\Database\NIST11.L
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Peak Number 5 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.58	25.16 ng	987355	Chrysene-d12	21.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	83
3		n-Decanoic acid	172	C10H20O2	000334-48-5	83
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	64
5		Undecanoic acid	186	C11H22O2	000112-37-8	64



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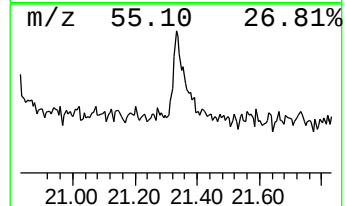
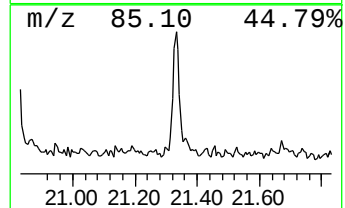
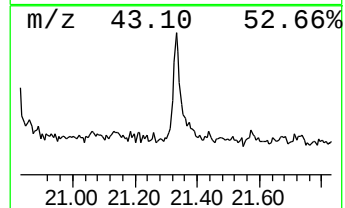
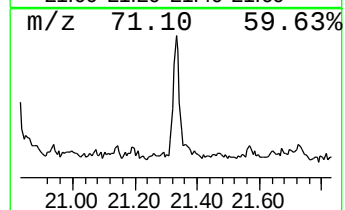
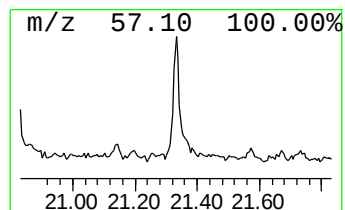
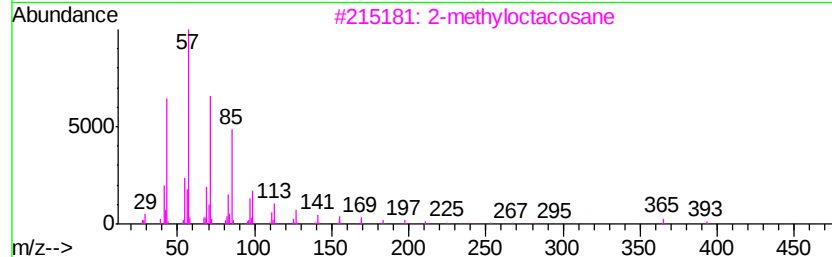
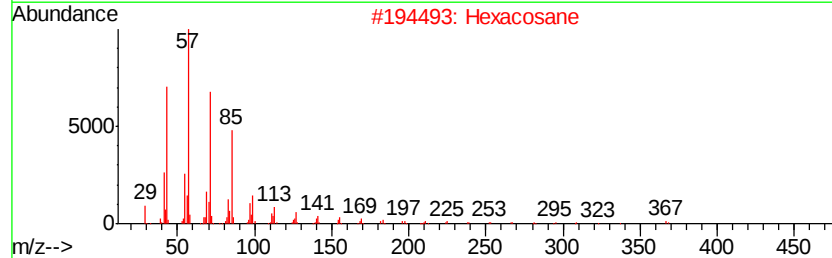
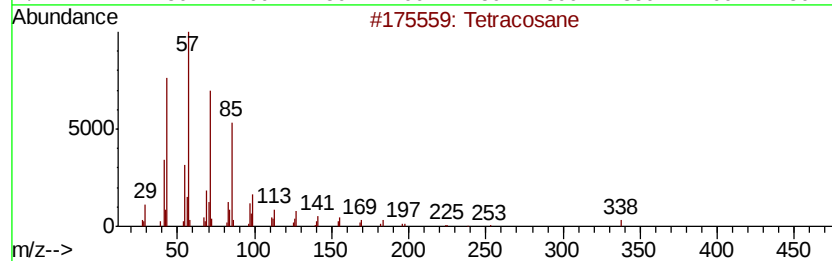
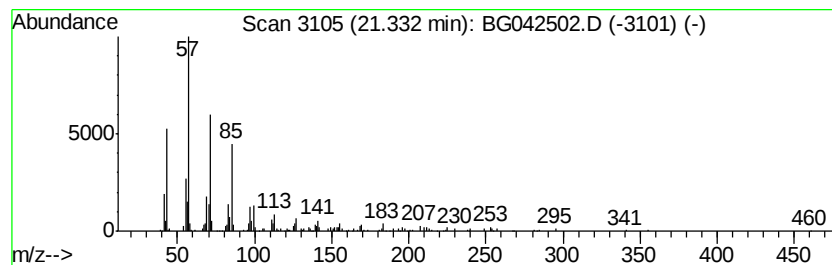
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TIC Library : C:\Database\NIST11.L
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Peak Number 6 Tetracosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.33	3.32 ng	130247	Chrysene-d12	21.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	93
2		Hexacosane	366	C26H54	000630-01-3	90
3		2-methyloctacosane	408	C29H60	1000376-72-8	90
4		Heneicosane, 11-decyl-	437	C31H64	055320-06-4	90
5		Octacosane	394	C28H58	000630-02-4	90



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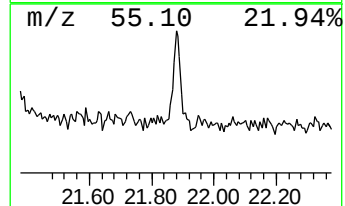
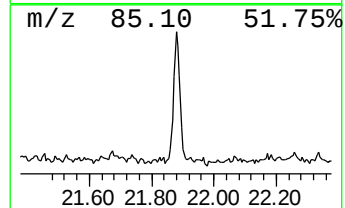
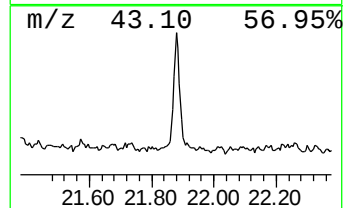
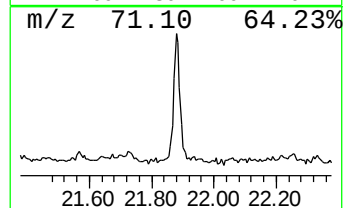
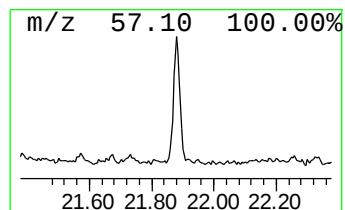
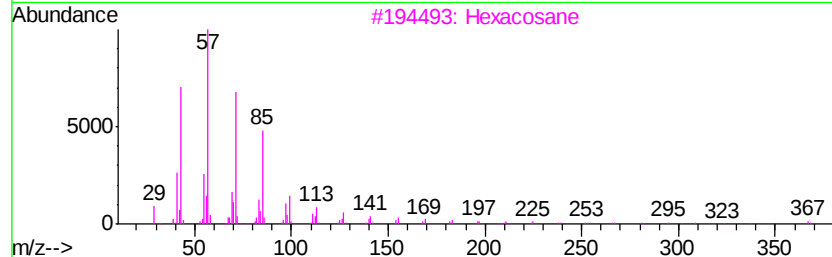
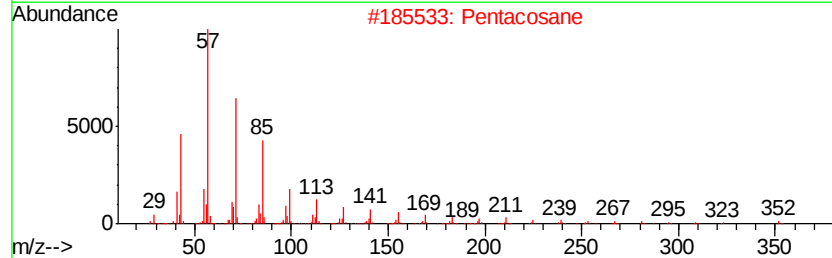
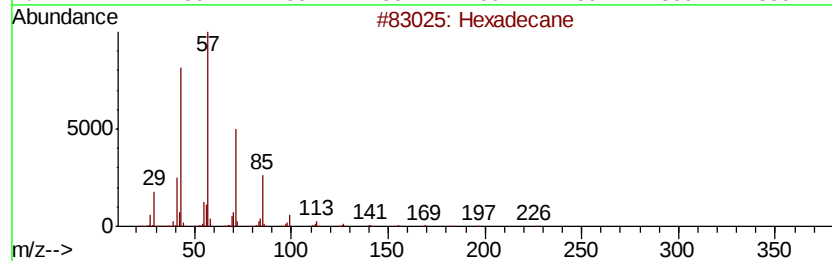
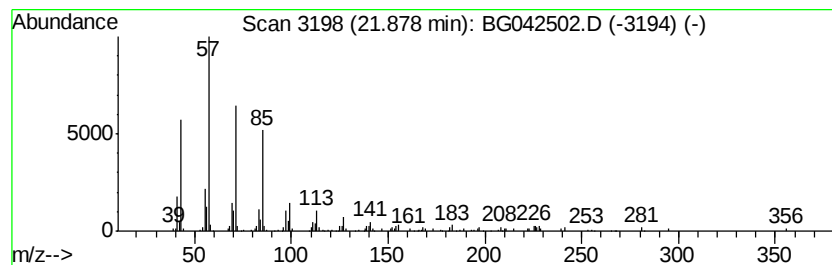
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.88	3.22 ng	126493	Chrysene-d12	21.68

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	95
2	Pentacosane		352	C25H52	000629-99-2	91
3	Hexacosane		366	C26H54	000630-01-3	90
4	Heptadecane		240	C17H36	000629-78-7	90
5	Tetracosane		338	C24H50	000646-31-1	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082619\
Data File : BG042502.D
Acq On : 26 Aug 2019 17:10
Operator : HP/JU
Sample : K4491-06
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
WP-1

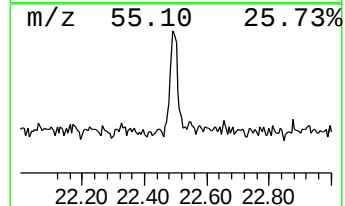
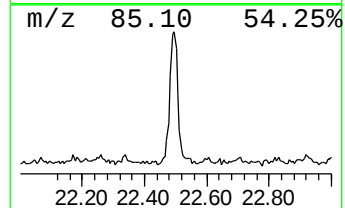
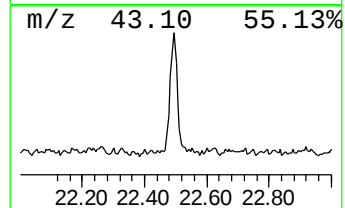
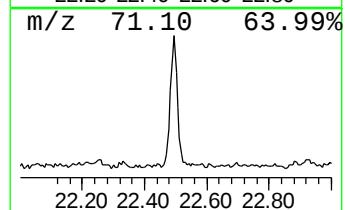
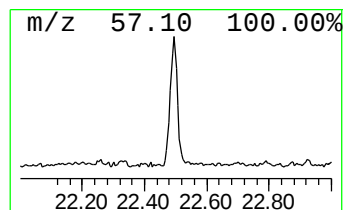
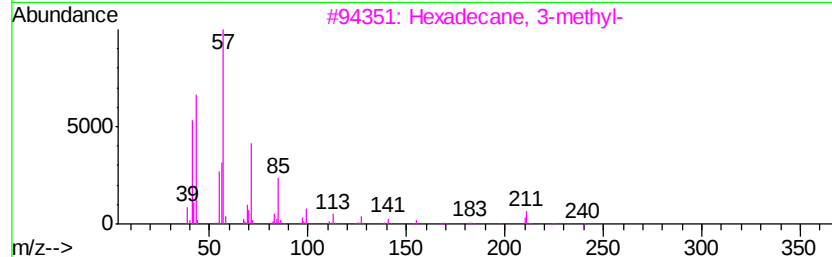
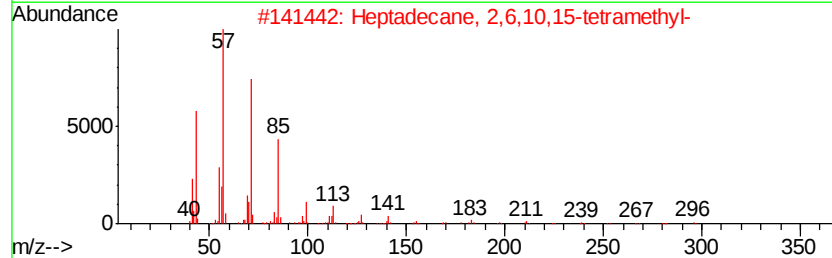
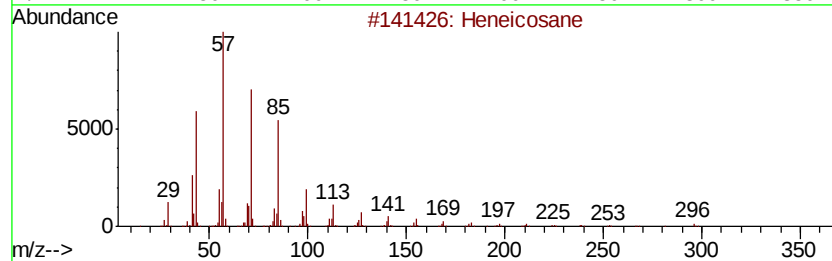
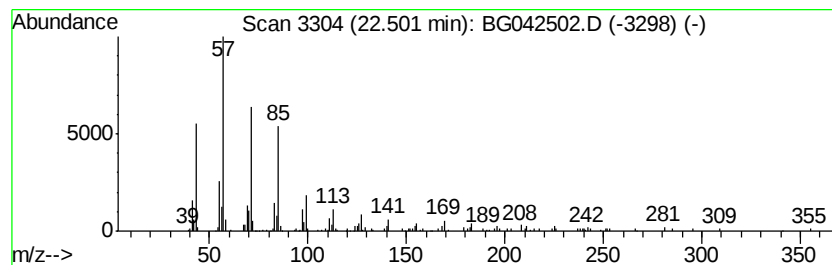
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heneicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.50	4.07 ng	159729	Chrysene-d12	21.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	98
2		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	93
3		Hexadecane, 3-methyl-	240	C17H36	006418-43-5	93
4		Nonadecane, 9-methyl-	282	C20H42	013287-24-6	93
5		triacontane	422	C30H62	000638-68-6	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082619\
Data File : BG042502.D
Acq On : 26 Aug 2019 17:10
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Instrument :
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ClientSampleId :
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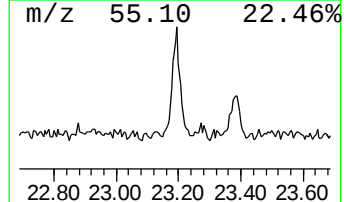
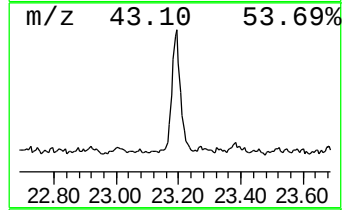
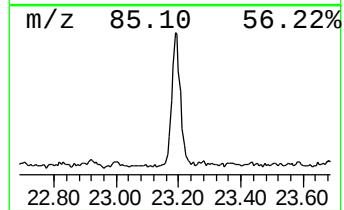
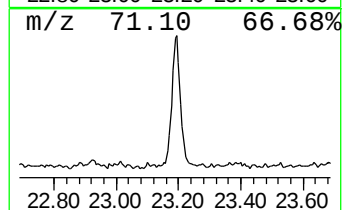
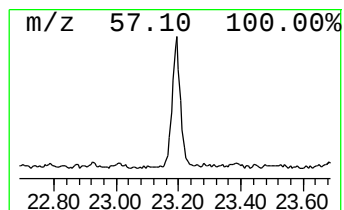
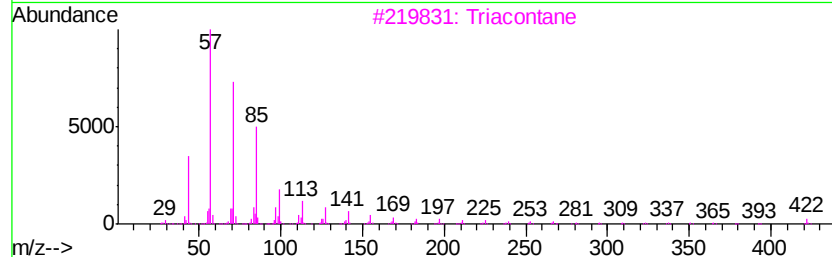
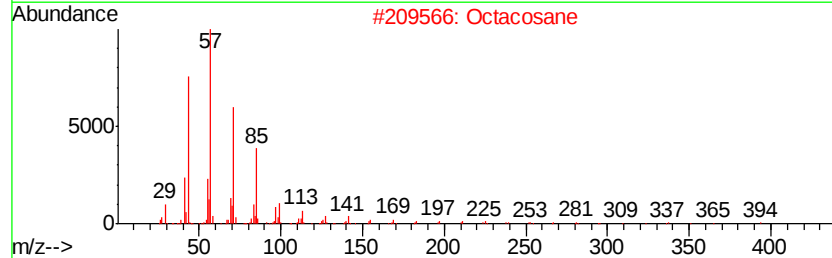
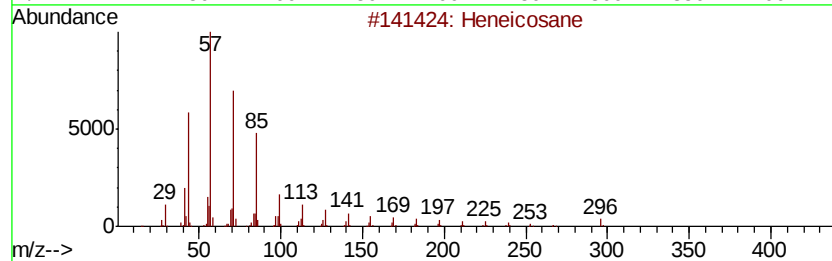
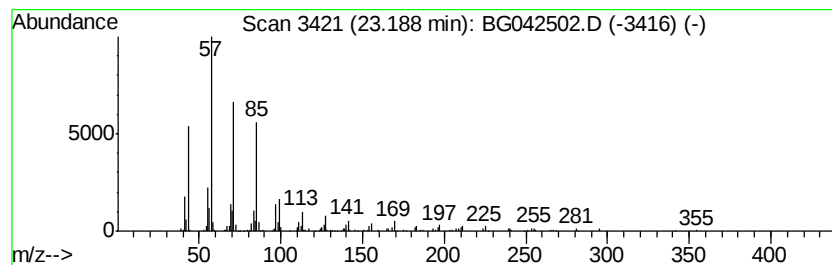
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.19	5.76 ng	226199	Chrysene-d12	21.68

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	96
2	Octacosane		394	C28H58	000630-02-4	91
3	Triacontane		422	C30H62	000638-68-6	91
4	Hentriacontane		437	C31H64	000630-04-6	91
5	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082619\
Data File : BG042502.D
Acq On : 26 Aug 2019 17:10
Operator : HP/JU
Sample : K4491-06
Misc :
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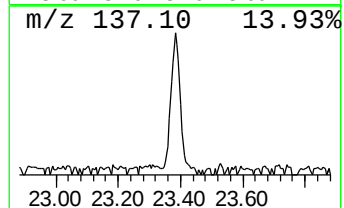
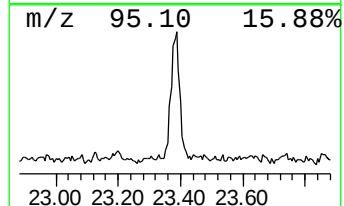
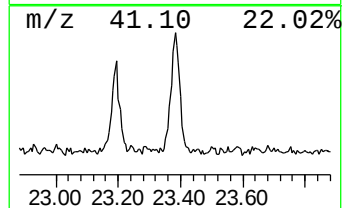
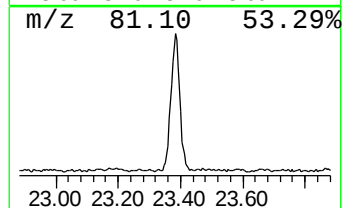
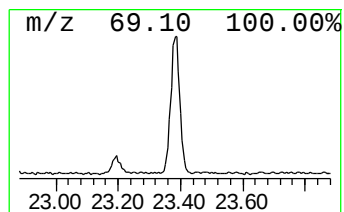
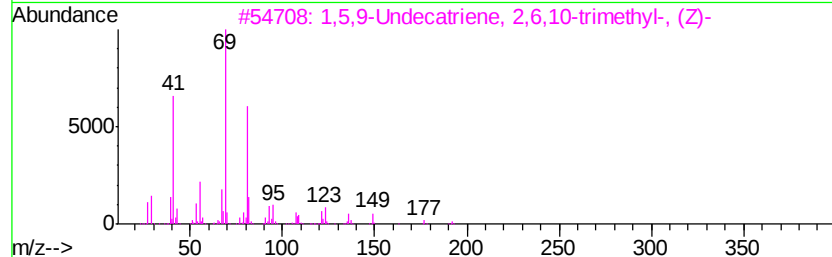
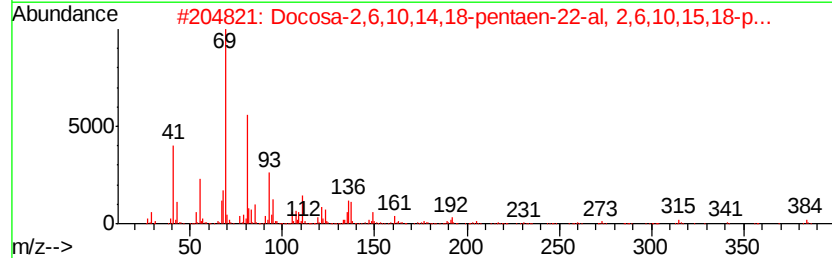
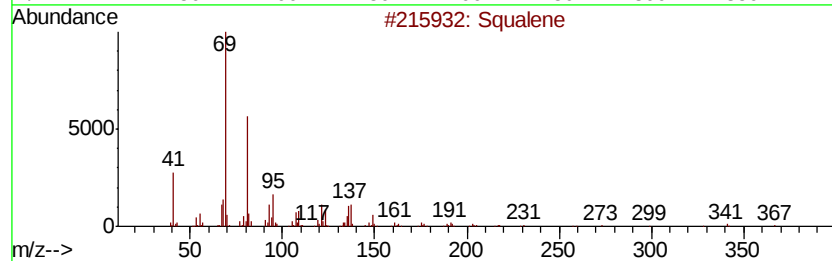
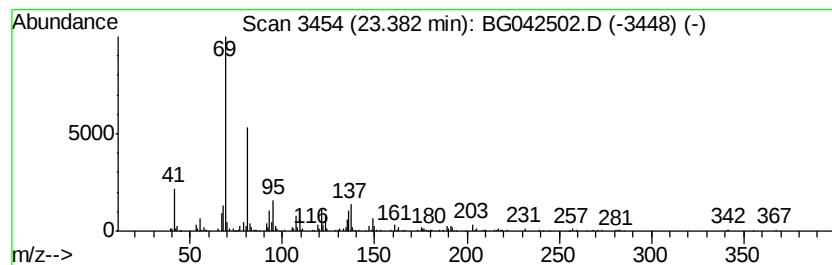
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Squalene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.38	4.04 ng	255677	Perylene-d12	24.93	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Squalene	410	C30H50	000111-02-4	91
2	Docosa-2,6,10,14,18-pentaen-22-a...	384	C27H44O	1000163-04-7	90
3	1,5,9-Undecatriene, 2,6,10-trime...	192	C14H24	062951-96-6	81
4	1,5,9-Decatriene, 2,3,5,8-tetram...	192	C14H24	230646-72-7	80
5	2,6,10,14-Hexadecatetraenoic aci...	332	C22H36O2	024035-35-6	78



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082619\
Data File : BG042502.D
Acq On : 26 Aug 2019 17:10
Operator : HP/JU
Sample : K4491-06
Misc :
ALS Vial : 14 Sample Multiplier: 1

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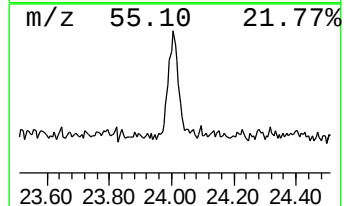
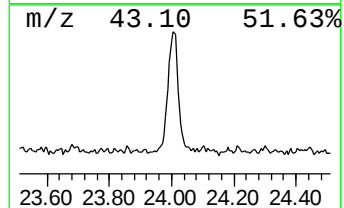
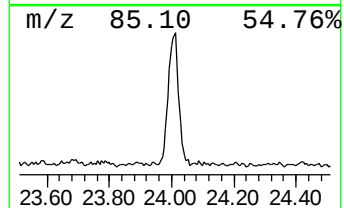
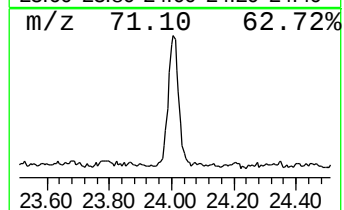
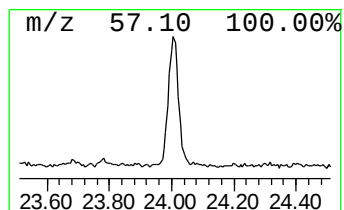
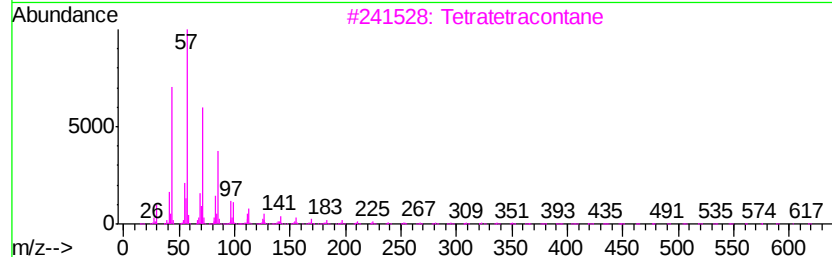
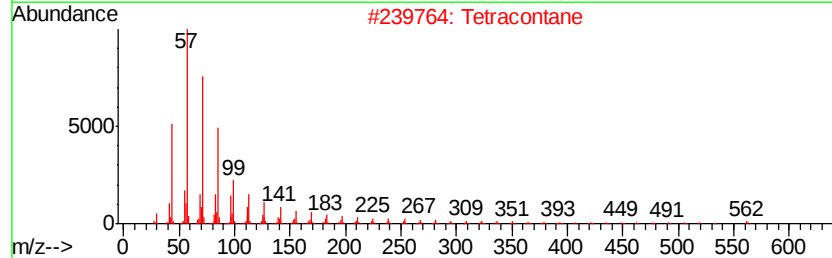
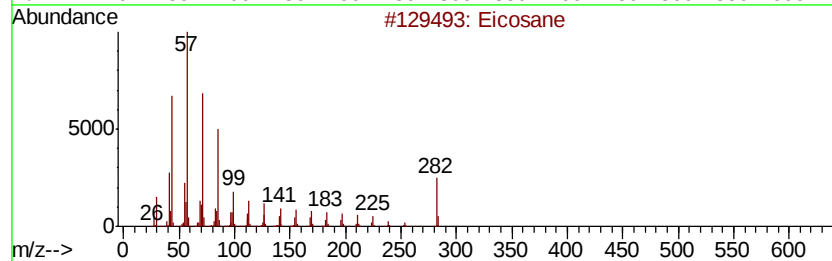
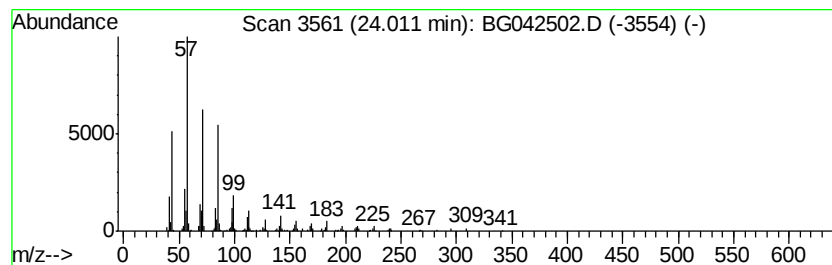
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Eicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.01	3.84 ng	242794	Perylene-d12	24.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	95
2		Tetracontane	563	C40H82	004181-95-7	91
3		Tetratetracontane	619	C44H90	007098-22-8	90
4		Hentriacontane	437	C31H64	000630-04-6	90
5		Pentacosane	352	C25H52	000629-99-2	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082619\
Data File : BG042502.D
Acq On : 26 Aug 2019 17:10
Operator : HP/JU
Sample : K4491-06
Misc :
ALS Vial : 14 Sample Multiplier: 1

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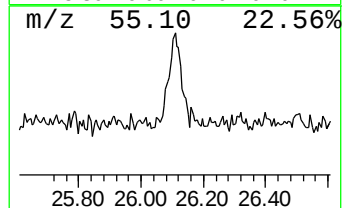
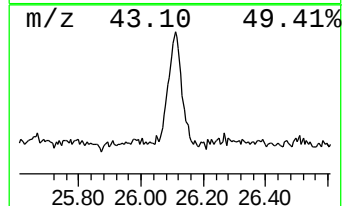
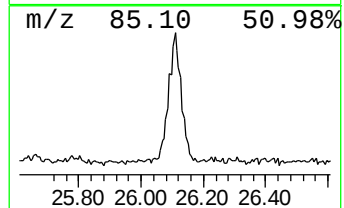
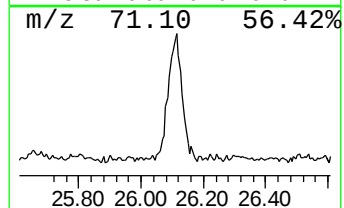
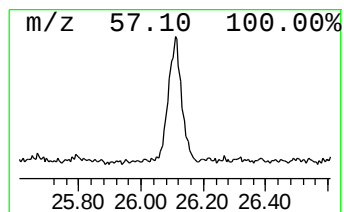
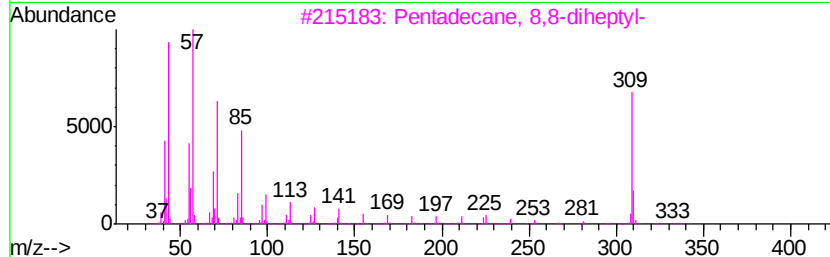
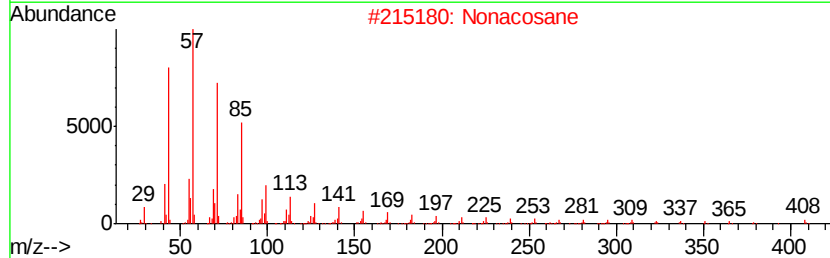
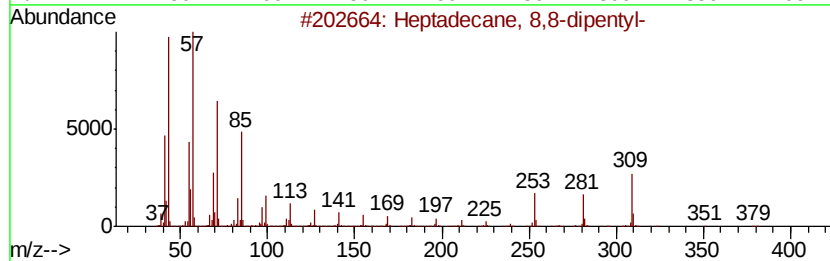
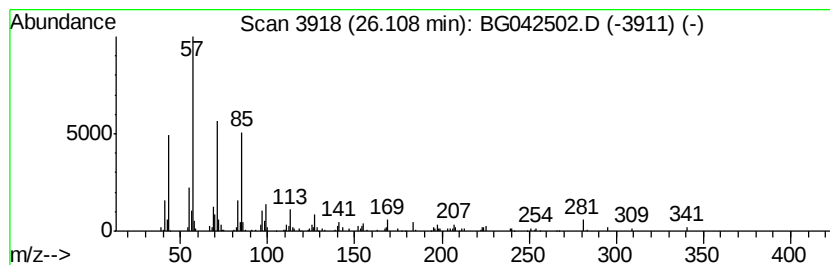
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heptadecane, 8,8-dipentyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.11	3.22 ng	203888	Perylene-d12	24.93

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 8,8-dipentyl-	380	C27H56	055282-28-5	90
2			Nonacosane	408	C29H60	000630-03-5	90
3			Pentadecane, 8,8-diheptyl-	408	C29H60	055268-72-9	86
4			2-methyloctacosane	408	C29H60	1000376-72-8	86
5			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG082619\
Data File : BG042502.D
Acq On : 26 Aug 2019 17:10
Operator : HP/JU
Sample : K4491-06
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG082219.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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unknown7.53	7.53	77.0	ng	745261	1	8.00	193688	20.0
n-Hexadecanoic acid	18.31	93.6	ng	2122080	4	17.41	453310	20.0
Octadecanoic acid	19.58	25.2	ng	987355	5	21.68	784947	20.0
Tetracosane	21.33	3.3	ng	130247	5	21.68	784947	20.0
Hexadecane	21.88	3.2	ng	126493	5	21.68	784947	20.0
Heneicosane	22.50	4.1	ng	159729	5	21.68	784947	20.0
Octacosane	23.19	5.8	ng	226199	5	21.68	784947	20.0
Squalene	23.38	4.0	ng	255677	6	24.93	1264770	20.0
Eicosane	24.01	3.8	ng	242794	6	24.93	1264770	20.0
Heptadecane, 8,8-...	26.11	3.2	ng	203888	6	24.93	1264770	20.0