

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010720\
 Data File : BG043957.D
 Acq On : 7 Jan 2020 16:41
 Operator : JU
 Sample : K6463-02 2X
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 GP-6-(4-6)

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	3.155	5	11	23	rVB	169759	313415	4.78%	0.512%
2	5.070	331	337	347	rVB	256143	394751	6.02%	0.645%
3	5.540	409	417	426	rBV	1081991	1744189	26.59%	2.848%
4	7.127	678	687	700	rBV	1146617	2031479	30.97%	3.317%
5	7.497	739	750	759	rBV	1105619	1935905	29.52%	3.161%
6	7.961	822	829	836	rBV	580801	972837	14.83%	1.589%
7	8.284	874	884	893	rBV	819851	1411783	21.53%	2.305%
8	9.112	1018	1025	1036	rVB	649813	1159426	17.68%	1.893%
9	10.763	1297	1306	1313	rBV	778687	1390427	21.20%	2.271%
10	13.214	1716	1723	1734	rBV	1481322	2254459	34.37%	3.682%
11	14.042	1857	1864	1869	rBV	103454	163360	2.49%	0.267%
12	14.588	1950	1957	1964	rBV2	1195253	1880490	28.67%	3.071%
13	16.075	2203	2210	2217	rBV	1264404	2036120	31.05%	3.325%
14	16.204	2219	2232	2234	rBV3	229703	755137	11.51%	1.233%
15	16.269	2238	2243	2248	rVV3	138986	379648	5.79%	0.620%
16	17.332	2417	2424	2428	rBV	1403970	2150619	32.79%	3.512%
17	17.373	2428	2431	2437	rVV	385124	542459	8.27%	0.886%
18	17.462	2438	2446	2452	rVB	83027	154965	2.36%	0.253%
19	18.208	2569	2573	2576	rBV2	52877	86773	1.32%	0.142%
20	18.255	2578	2581	2586	rVV	81867	119850	1.83%	0.196%
21	18.337	2588	2595	2601	rVV5	68674	178408	2.72%	0.291%
22	18.402	2602	2606	2610	rVB3	61149	106866	1.63%	0.175%
23	18.572	2625	2635	2648	rVB6	45081	201099	3.07%	0.328%
24	18.707	2654	2658	2661	rBV2	49863	69728	1.06%	0.114%
25	19.042	2702	2715	2733	rBV2	1531455	6558495	100.00%	10.710%
26	19.165	2733	2736	2743	rVB5	43677	79107	1.21%	0.129%
27	19.383	2763	2773	2788	rVB2	1581905	4039539	61.59%	6.597%
28	19.741	2825	2834	2842	rVB	689019	1172534	17.88%	1.915%
29	19.818	2844	2847	2855	rBV6	40353	98901	1.51%	0.162%
30	19.947	2864	2869	2874	rVB	2190138	2782221	42.42%	4.543%
31	20.323	2921	2933	2943	rBV2	2041350	6193228	94.43%	10.114%
32	21.316	3094	3102	3118	rVB	2121791	5747938	87.64%	9.387%
33	21.580	3139	3147	3152	rBV2	1455642	2758146	42.05%	4.504%
34	21.627	3152	3155	3160	rVB	206453	301414	4.60%	0.492%

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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

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35	21.798	3181	3184	3189	rVB	170890	251221	3.83%	0.410%
36	22.456	3282	3296	3306	rBV	1499785	4543363	69.27%	7.419%
37	23.078	3397	3402	3414	rVB	248736	490723	7.48%	0.801%
38	23.789	3517	3523	3532	rVB3	534340	1569030	23.92%	2.562%
39	24.712	3671	3680	3688	rBV2	825807	2215384	33.78%	3.618%

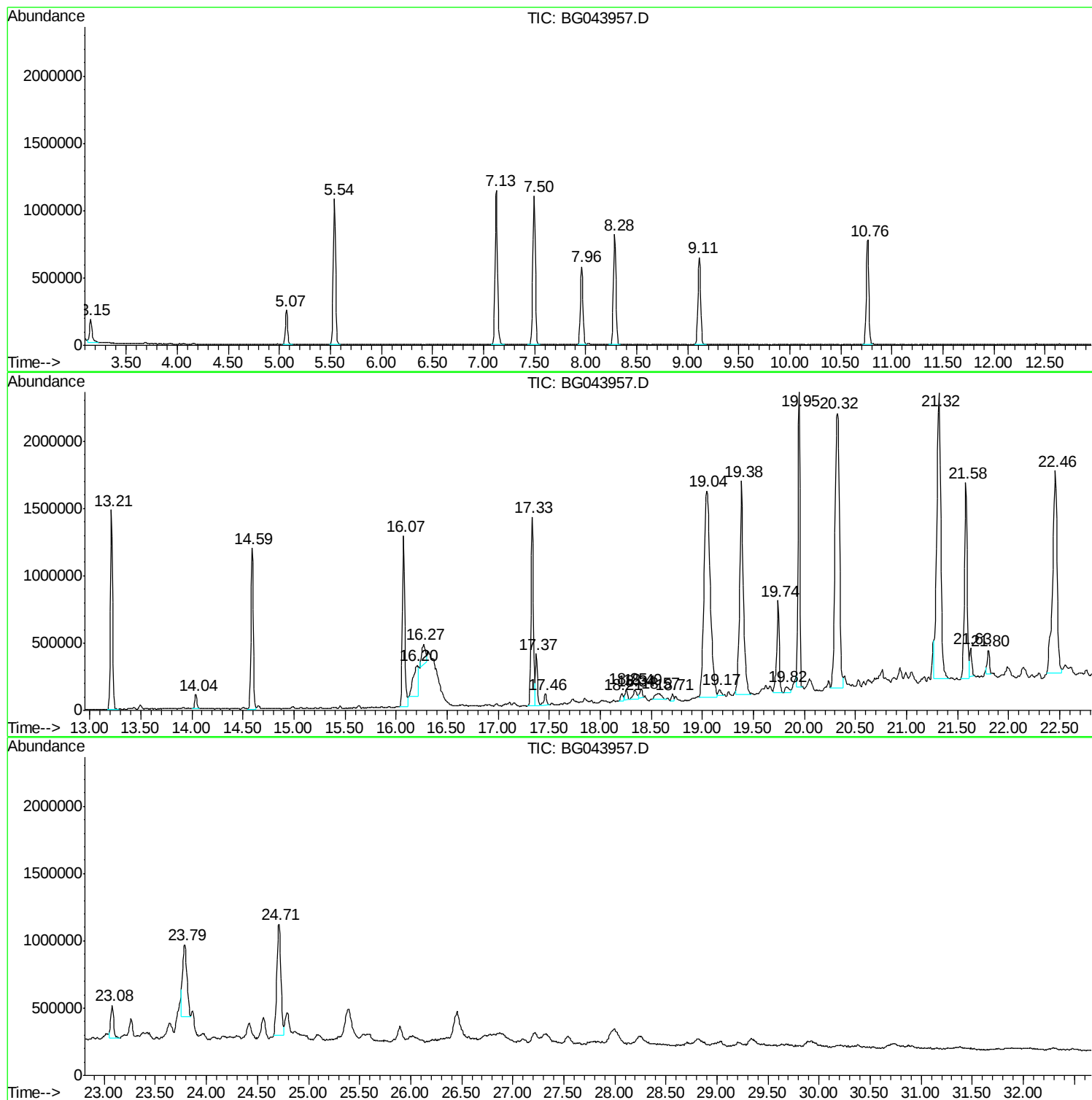
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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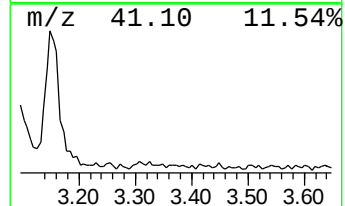
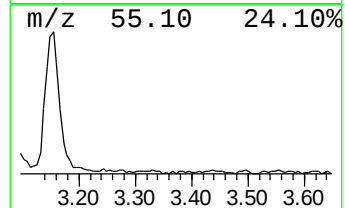
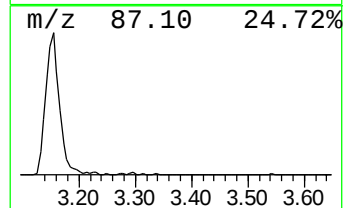
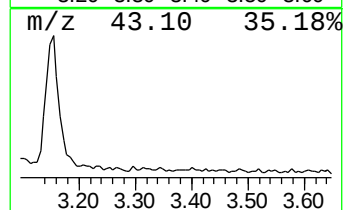
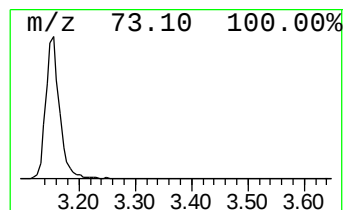
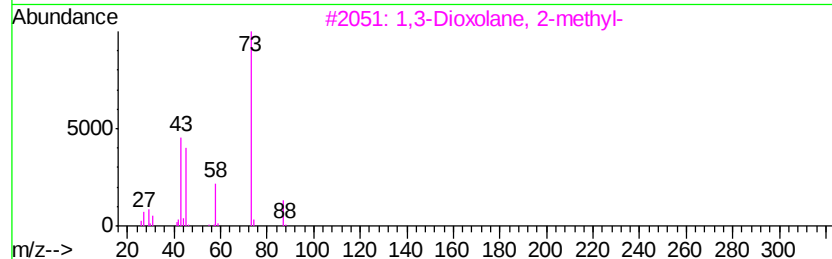
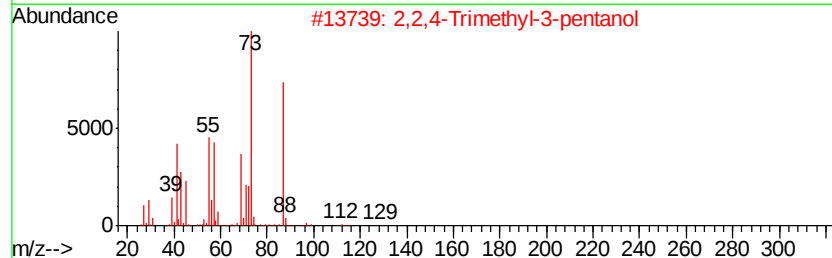
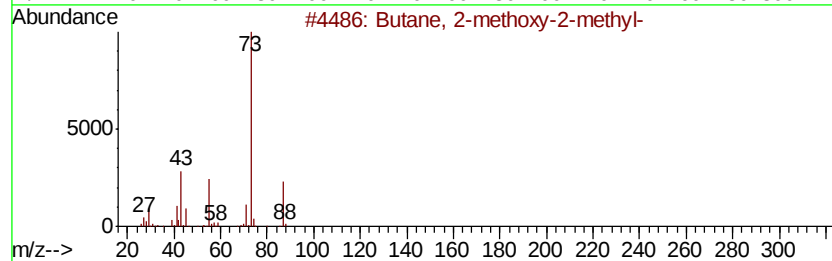
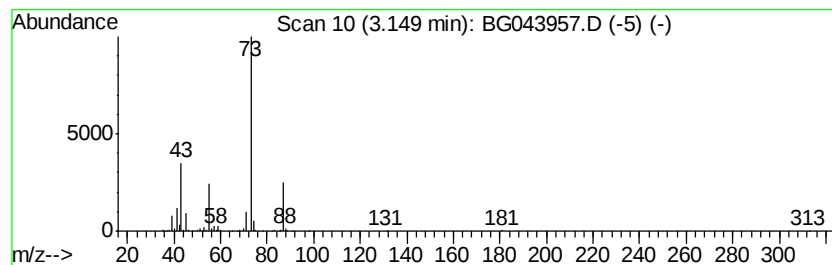
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.15	6.44 ng	313415	1,4-Dichlorobenzene-d4	7.96

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		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		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	12
4		N-Ethylformamide	73	C3H7NO	000627-45-2	9
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	9



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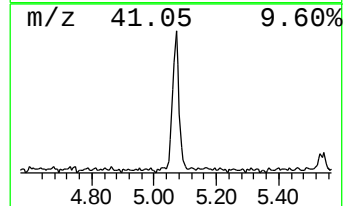
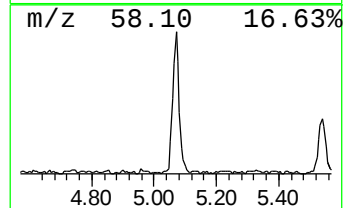
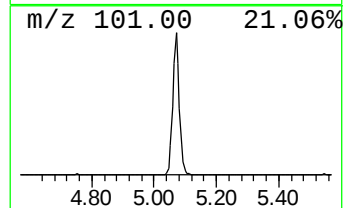
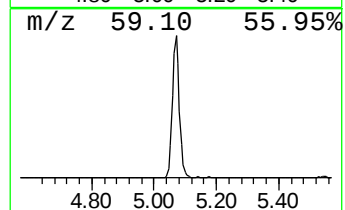
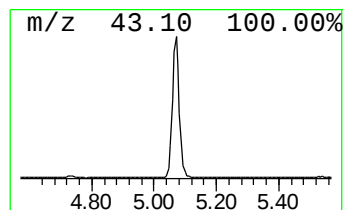
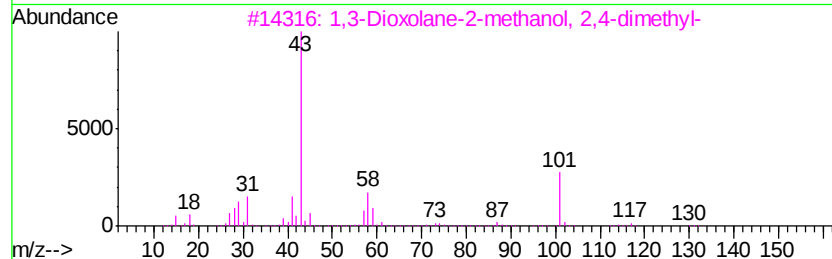
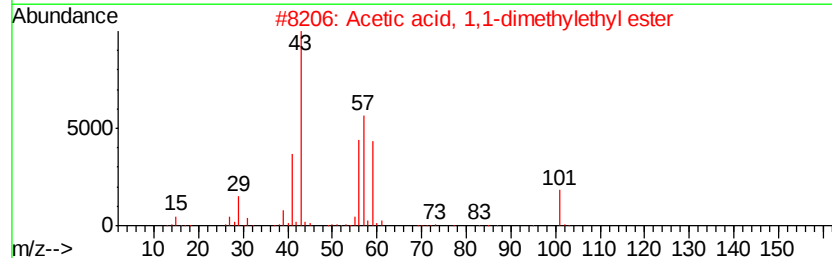
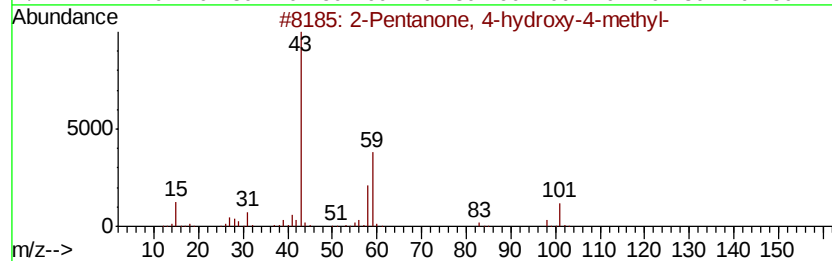
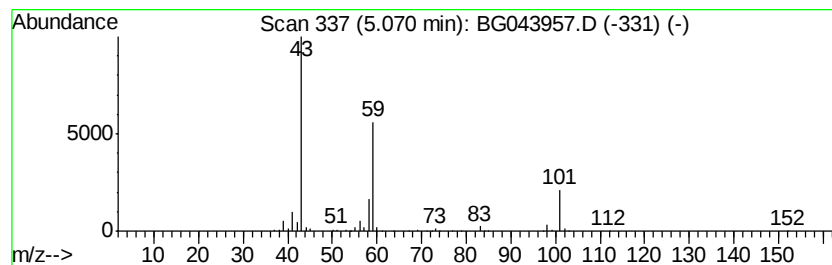
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	8.12 ng	394751	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	50
2			Acetic acid, 1,1-dimethylethyl ester	116	C6H12O2	000540-88-5	39
3			1,3-Dioxolane-2-methanol, 2,4-dimethyl-	132	C6H12O3	053951-43-2	28
4			Tetraolyme	222	C10H22O5	000143-24-8	25
5			4-Hydroxy-3-hexanone	116	C6H12O2	004984-85-4	23



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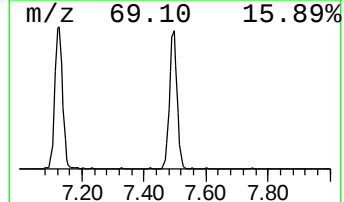
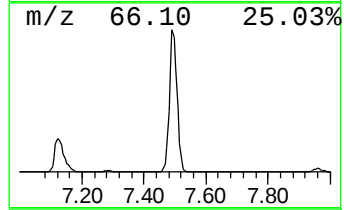
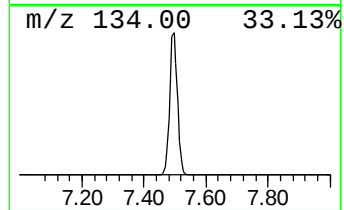
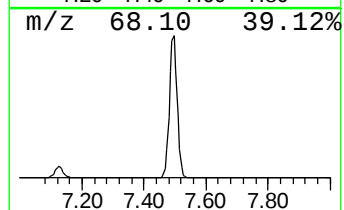
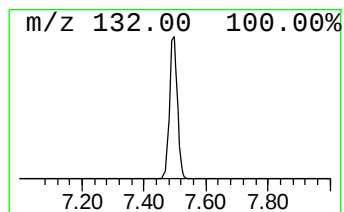
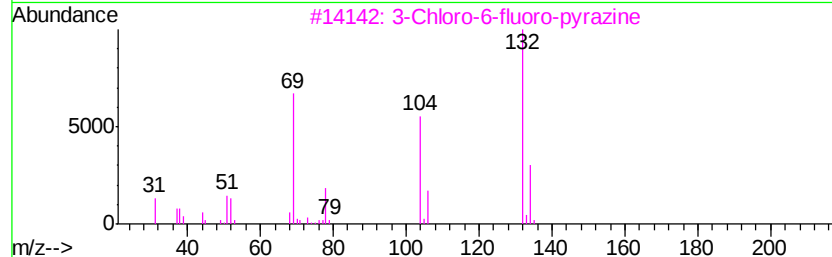
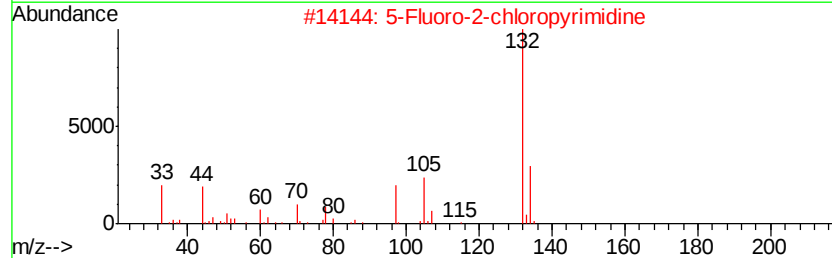
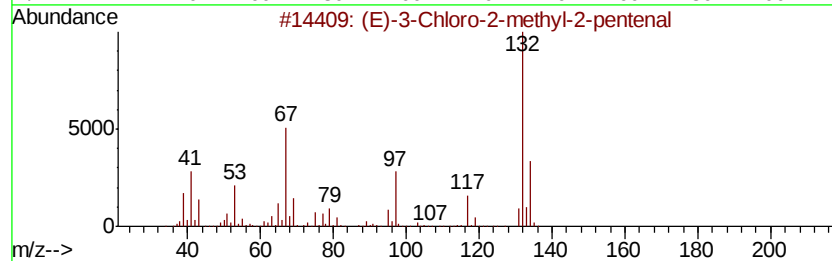
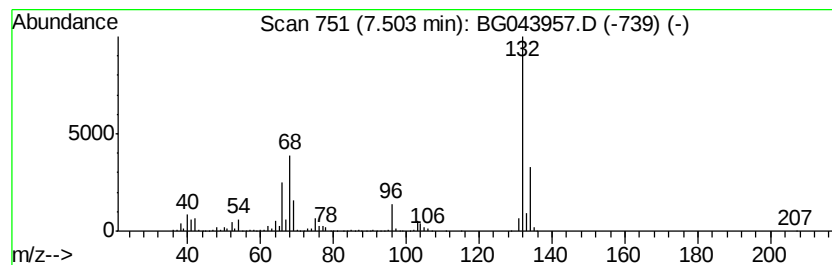
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 unknown7.50 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.50	39.80 ng	1935910	1,4-Dichlorobenzene-d4	7.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	30
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
4		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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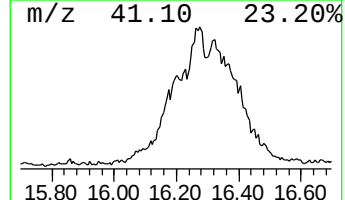
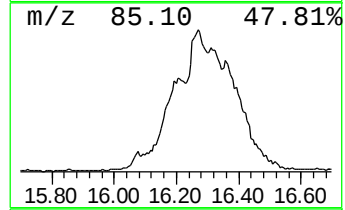
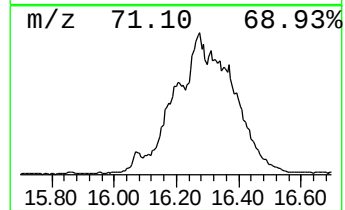
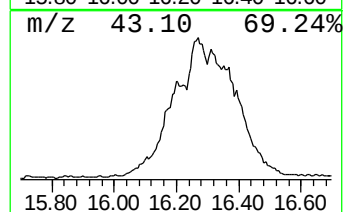
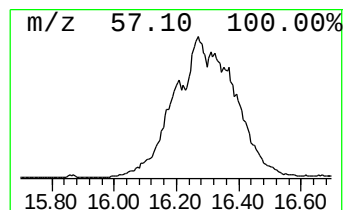
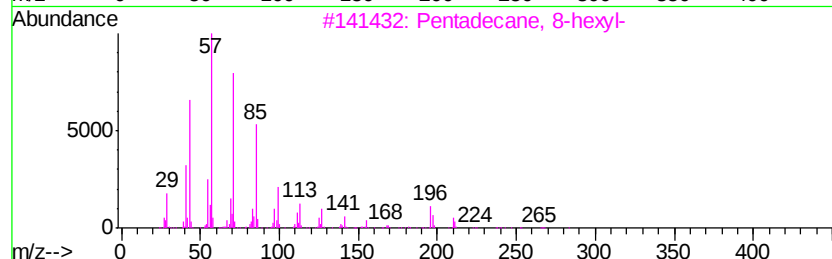
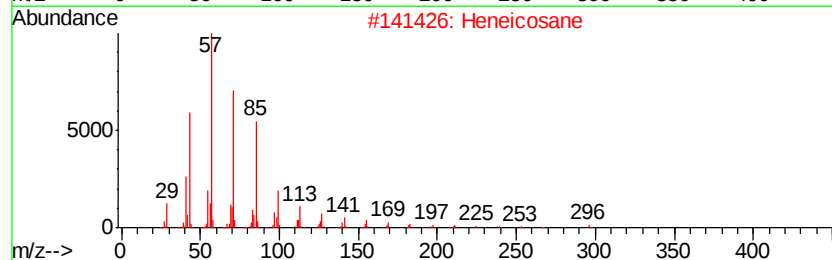
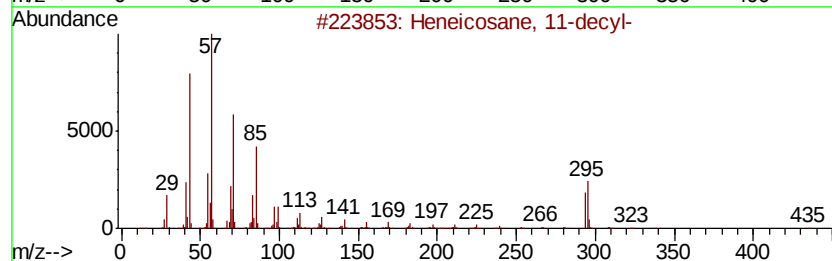
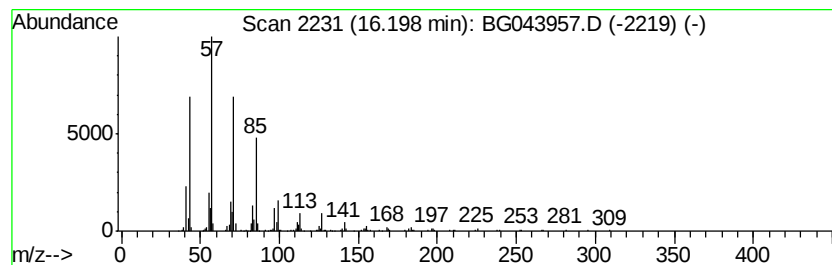
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 Peak Number 5 Heneicosane, 11-decyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
16.20	7.02 ng	755137	Phenanthrene-d10		17.33

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane, 11-decyl-	437	C31H64	055320-06-4	91
2			Heneicosane	296	C21H44	000629-94-7	91
3			Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	91
4			Pentadecane	212	C15H32	000629-62-9	90
5			Octadecane	254	C18H38	000593-45-3	87



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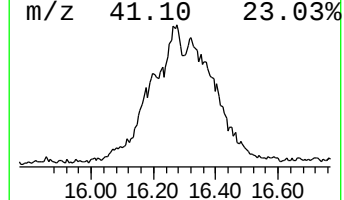
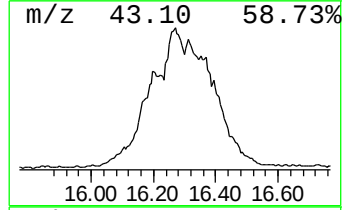
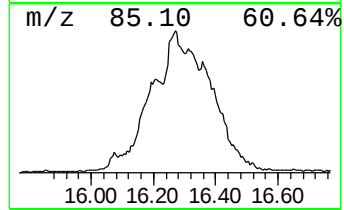
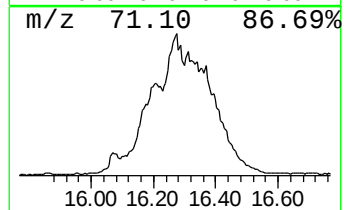
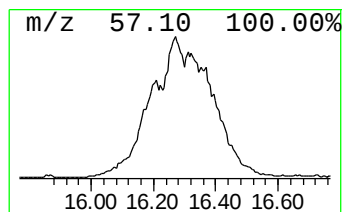
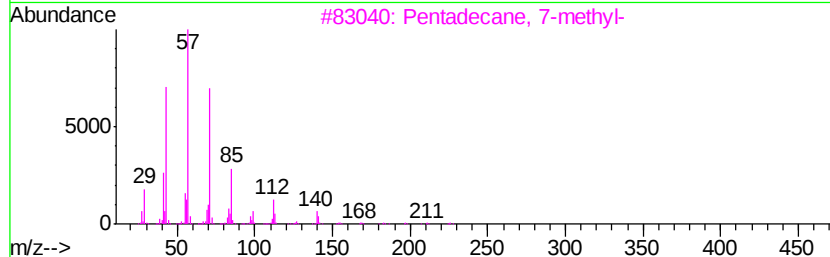
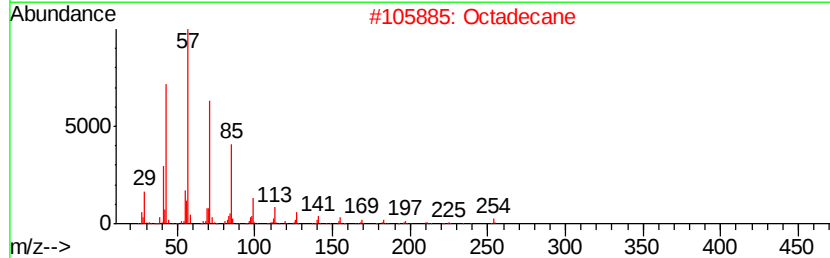
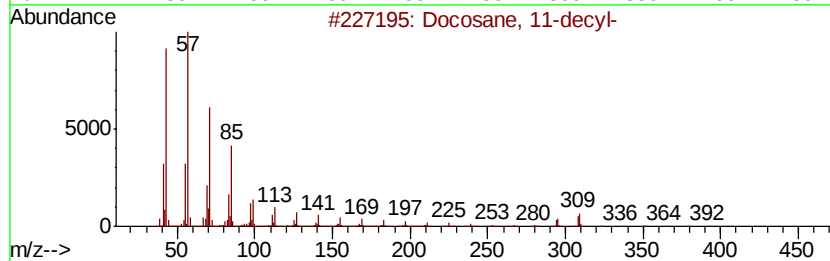
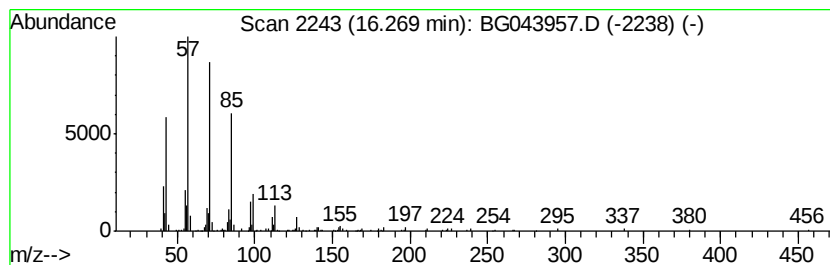
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Peak Number 6 Docosane, 11-decyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.		
16.27	3.53 ng	379648	Phenanthrene-d10		17.33		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Docosane, 11-decyl-	451	C32H66	055401-55-3	90
2			Octadecane	254	C18H38	000593-45-3	86
3			Pentadecane, 7-methyl-	226	C16H34	006165-40-8	62
4			Heptadecane	240	C17H36	000629-78-7	58
5			Heptadecane, 7-methyl-	254	C18H38	020959-33-5	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010720\
Data File : BG043957.D
Acq On : 7 Jan 2020 16:41
Operator : JU
Sample : K6463-02 2X
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GP-6-(4-6)

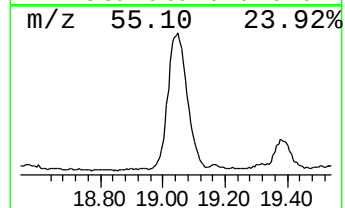
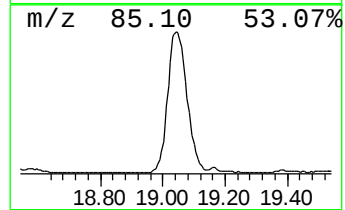
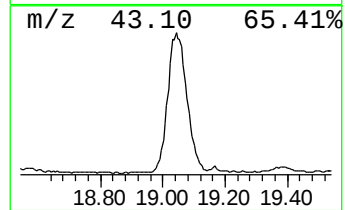
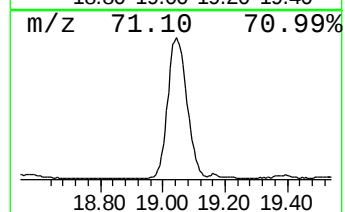
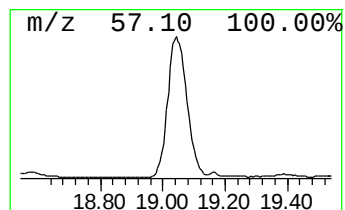
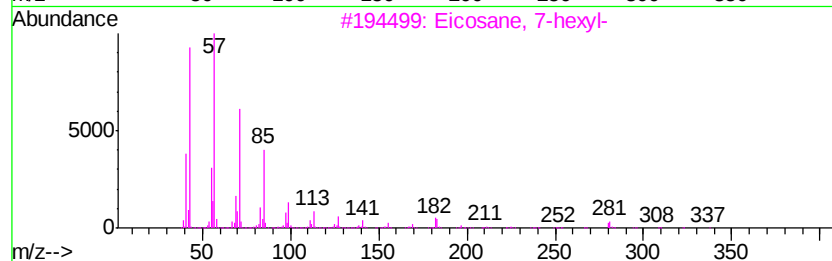
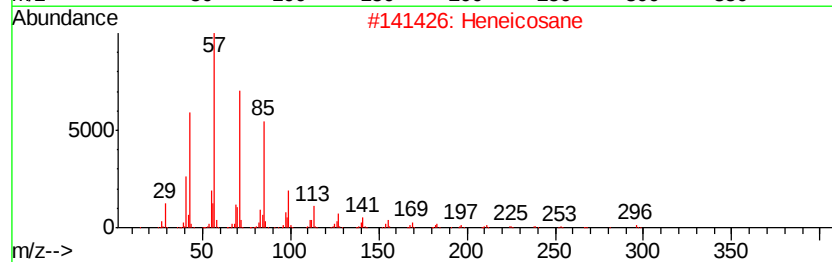
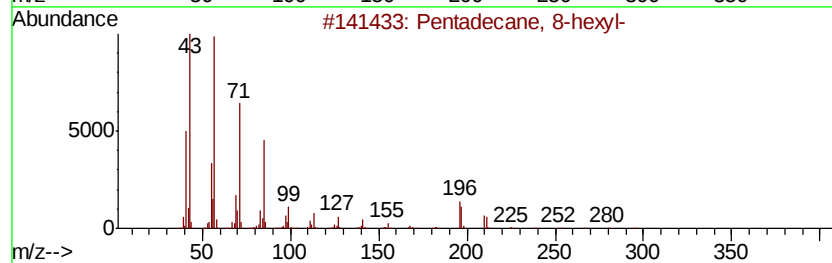
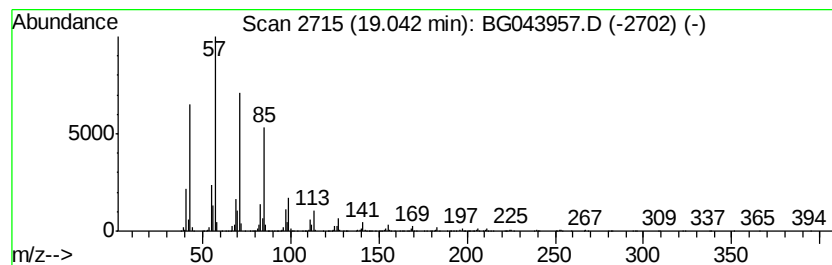
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Pentadecane, 8-hexyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.04	60.99 ng	6558500	Phenanthrene-d10	17.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
2		Heneicosane	296	C21H44	000629-94-7	91
3		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010720\
Data File : BG043957.D
Acq On : 7 Jan 2020 16:41
Operator : JU
Sample : K6463-02 2X
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GP-6-(4-6)

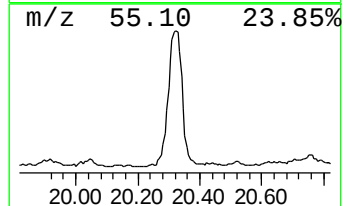
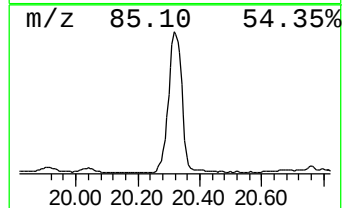
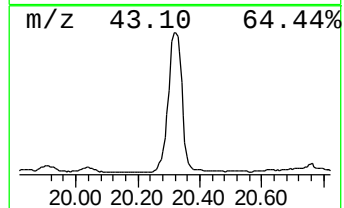
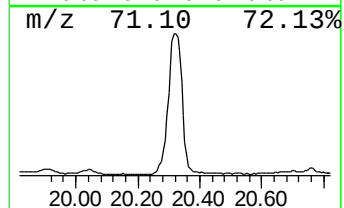
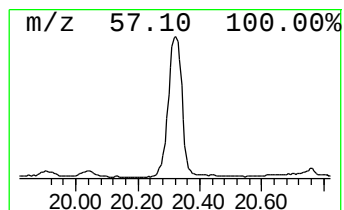
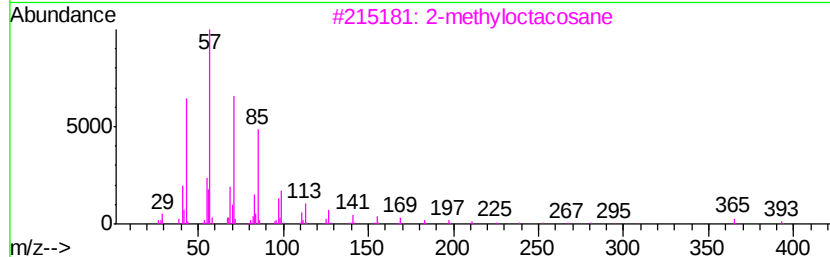
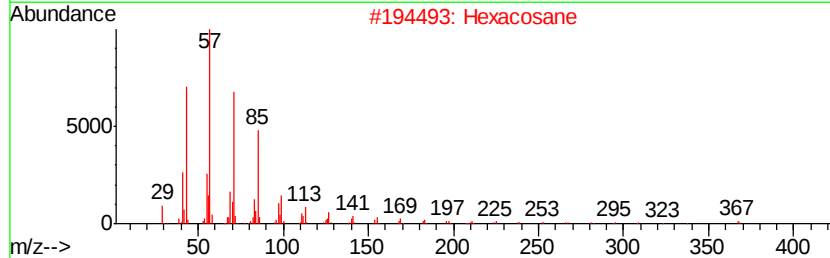
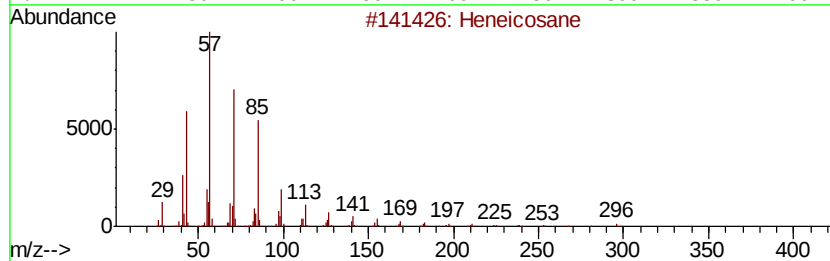
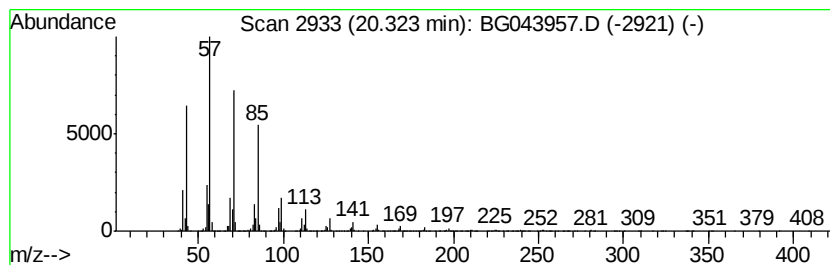
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexacosane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.32	44.91 ng	6193230	Chrysene-d12	21.58

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	Heneicosane, 11-(1-ethylpropyl)-		366	C26H54	055282-11-6	91
5	Hexatriacontane		507	C36H74	000630-06-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010720\
 Data File : BG043957.D
 Acq On : 7 Jan 2020 16:41
 Operator : JU
 Sample : K6463-02 2X
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

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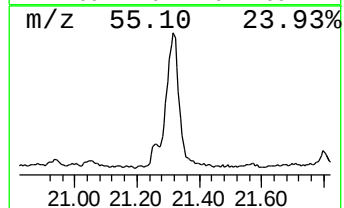
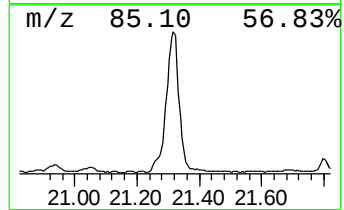
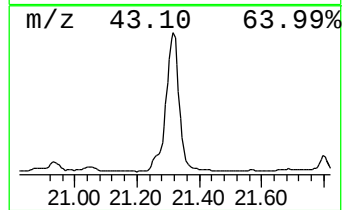
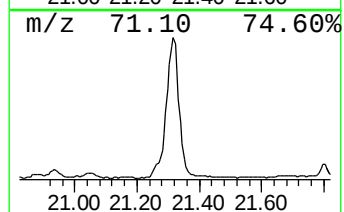
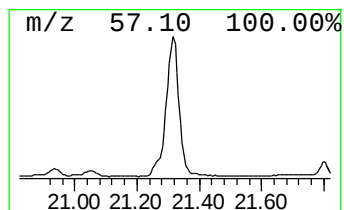
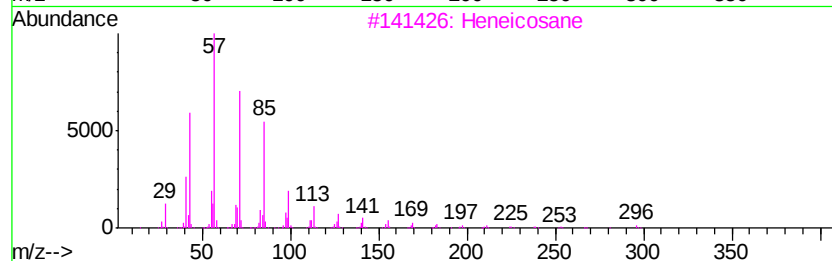
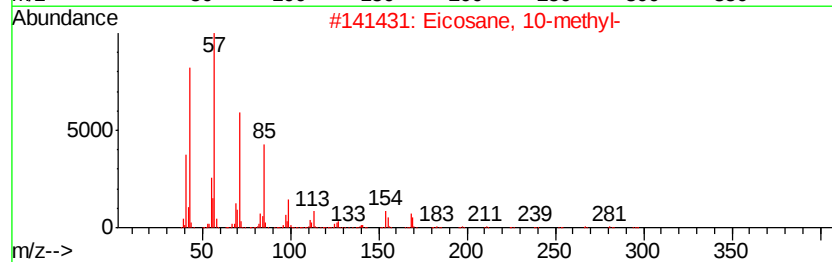
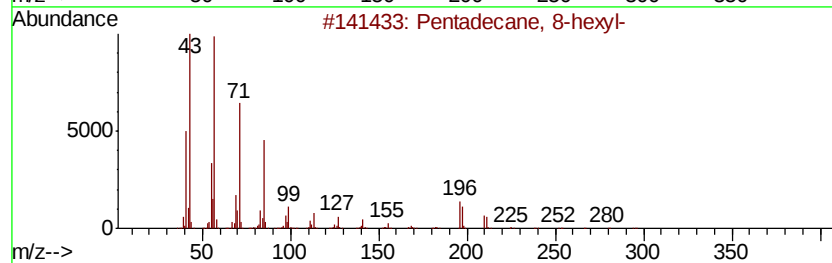
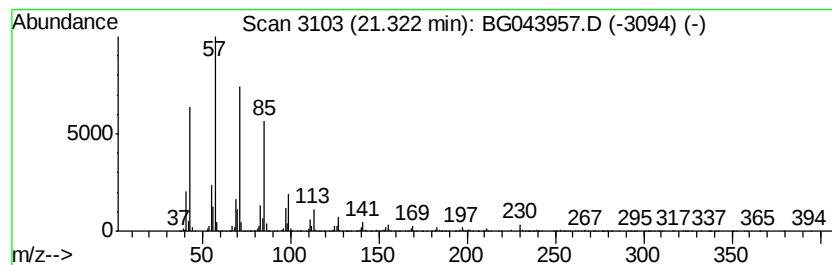
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 9 Eicosane, 10-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.32	41.68 ng	5747940	Chrysene-d12	21.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
2		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
3		Heneicosane	296	C21H44	000629-94-7	91
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010720\
Data File : BG043957.D
Acq On : 7 Jan 2020 16:41
Operator : JU
Sample : K6463-02 2X
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
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ClientSampleId :
GP-6-(4-6)

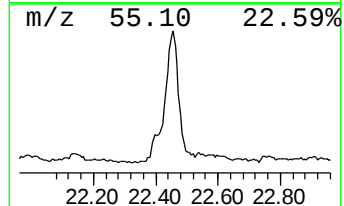
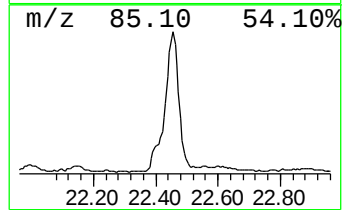
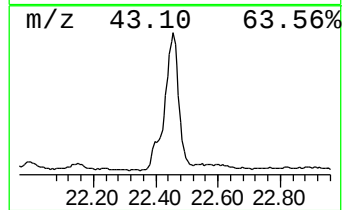
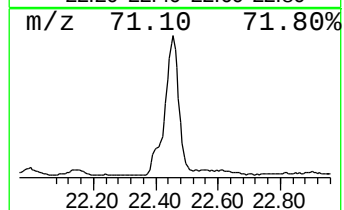
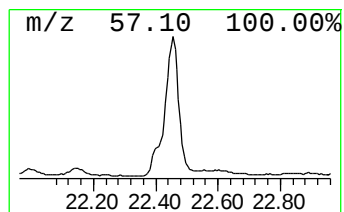
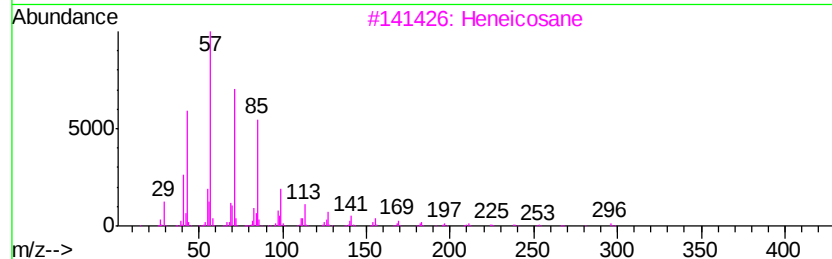
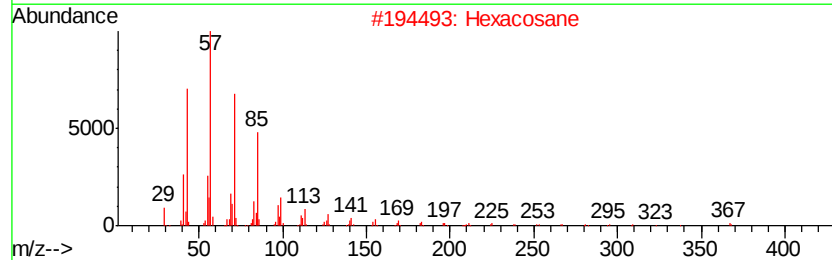
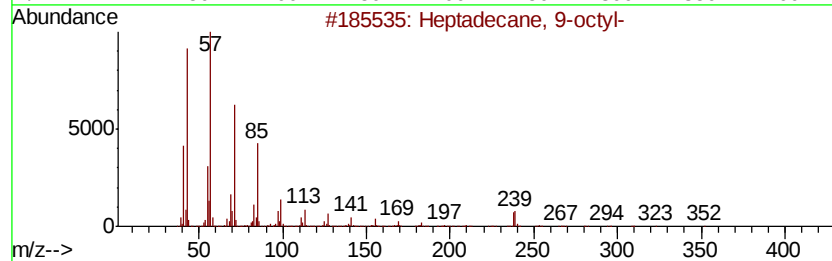
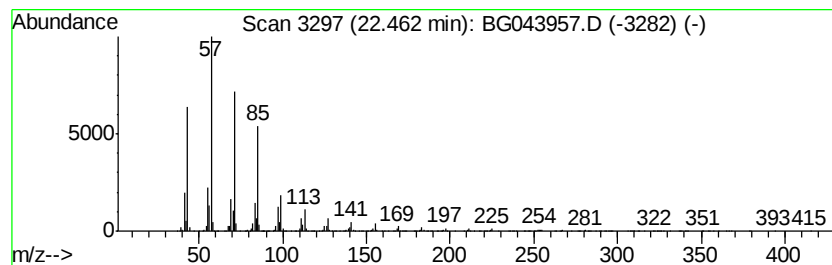
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptadecane, 9-octyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.46	32.95 ng	4543360	Chrysene-d12	21.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
2		Hexacosane	366	C26H54	000630-01-3	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		Tetratetracontane	619	C44H90	007098-22-8	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010720\
Data File : BG043957.D
Acq On : 7 Jan 2020 16:41
Operator : JU
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Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
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ClientSampleId :
GP-6-(4-6)

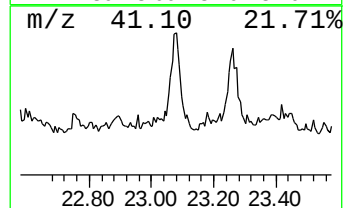
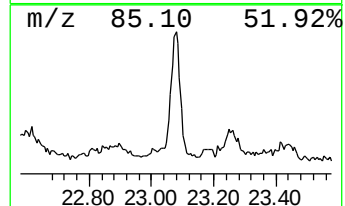
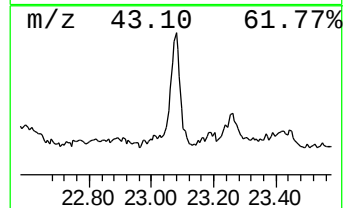
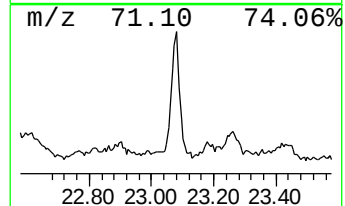
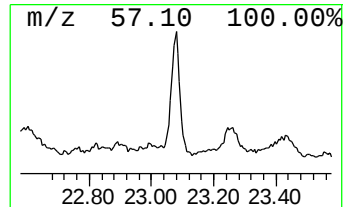
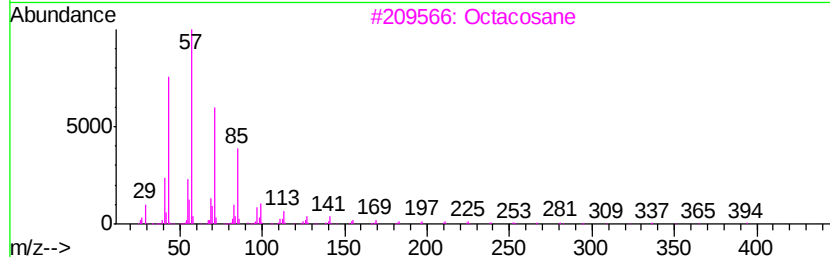
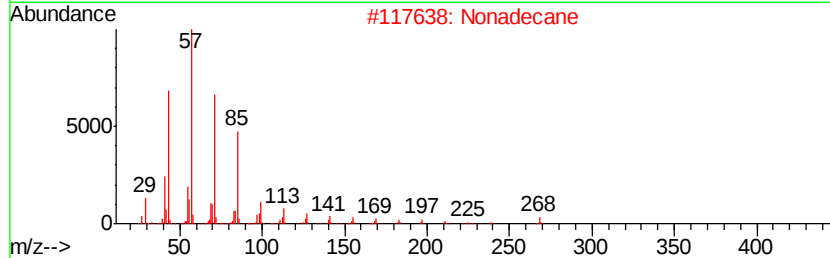
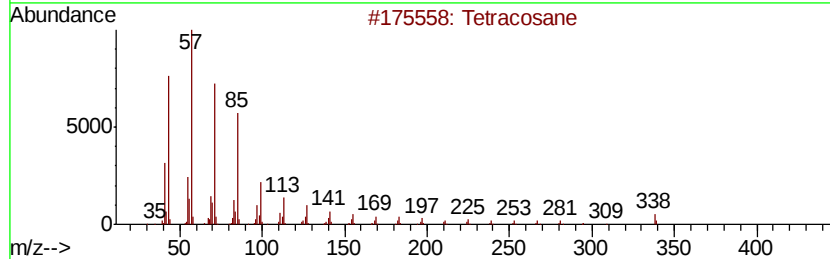
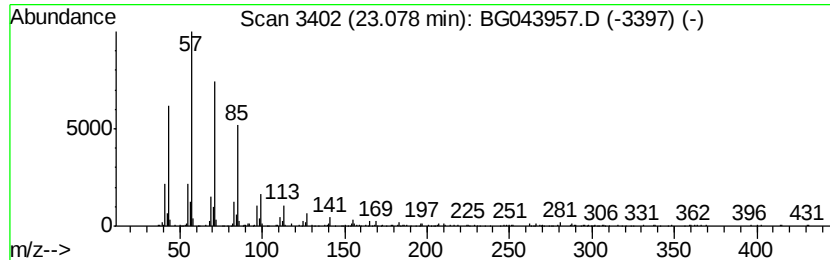
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tetracosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.08	3.56 ng	490723	Chrysene-d12	21.58

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	97
2			Nonadecane	268	C19H40	000629-92-5	93
3			Octacosane	394	C28H58	000630-02-4	91
4			Hexacosane	366	C26H54	000630-01-3	91
5			Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010720\
Data File : BG043957.D
Acq On : 7 Jan 2020 16:41
Operator : JU
Sample : K6463-02 2X
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
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ClientSampleId :
GP-6-(4-6)

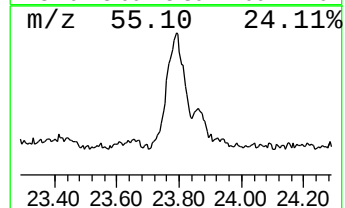
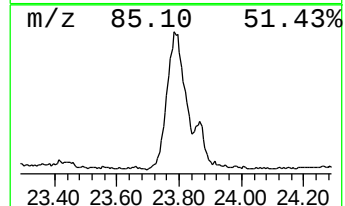
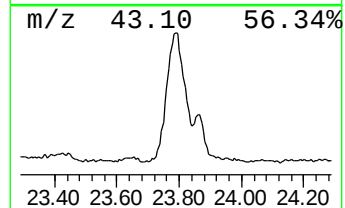
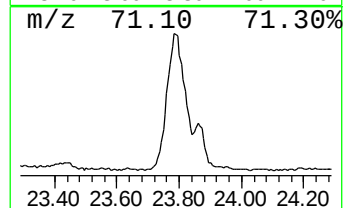
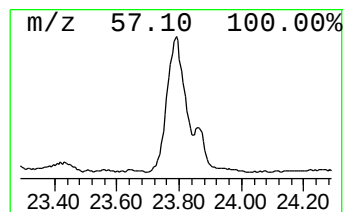
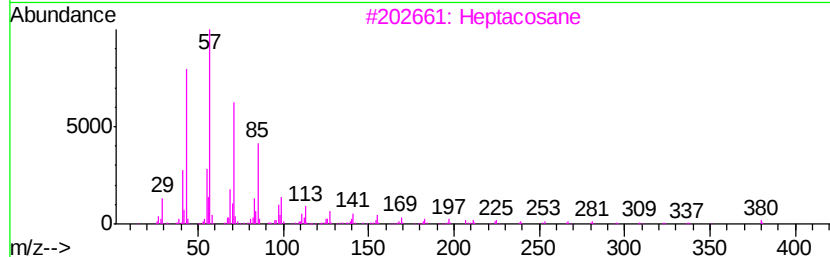
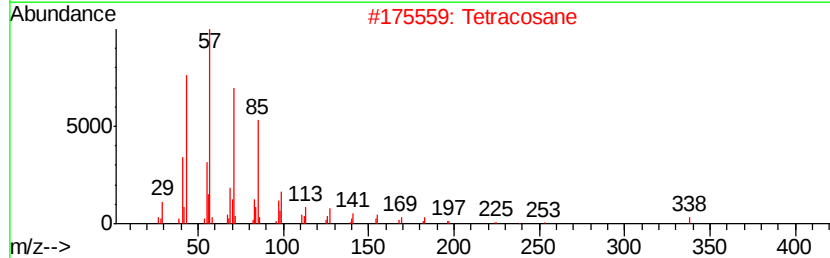
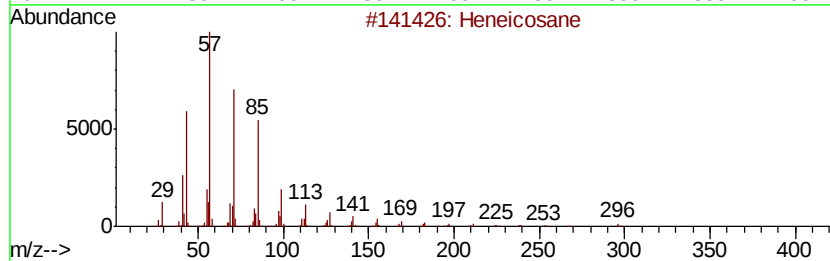
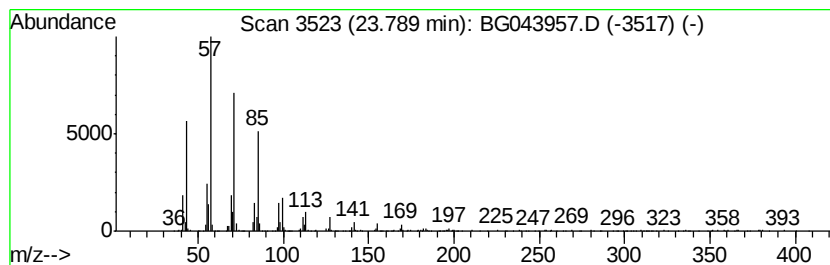
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heneicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.79	14.16 ng	1569030	Perylene-d12	24.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	97
2		Tetracosane	338	C24H50	000646-31-1	95
3		Heptacosane	380	C27H56	000593-49-7	91
4		Hexacosane, 9-octyl-	479	C34H70	055429-83-9	91
5		Docosane, 11-decyl-	451	C32H66	055401-55-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG010720\
 Data File : BG043957.D
 Acq On : 7 Jan 2020 16:41
 Operator : JU
 Sample : K6463-02 2X
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 GP-6-(4-6)

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	8.1	ng	394751	1	7.96	972837	20.0
unknown7.50	7.50	39.8	ng	1935910	1	7.96	972837	20.0
Heneicosane, 11-d...	16.20	7.0	ng	755137	4	17.33	2150620	20.0
Docosane, 11-decyl-	16.27	3.5	ng	379648	4	17.33	2150620	20.0
Pentadecane, 8-he...	19.04	61.0	ng	6558500	4	17.33	2150620	20.0
Hexacosane	20.32	44.9	ng	6193230	5	21.58	2758150	20.0
Eicosane, 10-methyl-	21.32	41.7	ng	5747940	5	21.58	2758150	20.0
Heptadecane, 9-oc...	22.46	33.0	ng	4543360	5	21.58	2758150	20.0
Tetracosane	23.08	3.6	ng	490723	5	21.58	2758150	20.0
Heneicosane	23.79	14.2	ng	1569030	6	24.71	2215380	20.0