

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050919\
 Data File : BG040827.D
 Acq On : 9 May 2019 22:18
 Operator : HP/JU
 Sample : K2762-16
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 P021-SS003-0612-01

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-BG042319.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	3.152	6	11	19	rBV	135913	231526	5.13%	0.802%
2	3.229	19	24	35	rVB	236300	322532	7.15%	1.118%
3	5.673	432	440	449	rBV	1116302	1748264	38.76%	6.058%
4	6.795	624	631	639	rVB	142557	236063	5.23%	0.818%
5	7.277	705	713	729	rBV	1219844	2173642	48.19%	7.532%
6	7.565	755	762	771	rVB2	106772	178819	3.96%	0.620%
7	7.665	771	779	788	rBV	1142862	1937778	42.96%	6.714%
8	8.140	853	860	871	rBV	199429	348628	7.73%	1.208%
9	8.464	908	915	924	rVB	789994	1333228	29.56%	4.620%
10	9.316	1051	1060	1071	rBV	738372	1351353	29.96%	4.682%
11	10.967	1333	1341	1349	rBV	274871	500173	11.09%	1.733%
12	13.393	1746	1754	1771	rBV	1916600	2912401	64.56%	10.092%
13	14.210	1888	1893	1900	rBV	157328	214905	4.76%	0.745%
14	14.768	1981	1988	1999	rBV2	517731	832557	18.46%	2.885%
15	16.254	2230	2241	2254	rBV	1865676	2869385	63.61%	9.942%
16	17.512	2449	2455	2466	rBV2	743879	1125729	24.96%	3.901%
17	18.370	2594	2601	2607	rBV	269151	412494	9.14%	1.429%
18	19.627	2811	2815	2824	rBV	154433	261170	5.79%	0.905%
19	20.109	2890	2897	2909	rBV	3397854	4510850	100.00%	15.630%
20	20.414	2939	2949	2959	rBV	60501	157628	3.49%	0.546%
21	20.714	2996	3000	3006	rBV9	47716	77066	1.71%	0.267%
22	20.855	3015	3024	3033	rBV2	352953	632757	14.03%	2.193%
23	21.254	3087	3092	3100	rVV4	193024	297633	6.60%	1.031%
24	21.390	3111	3115	3119	rBV4	50556	69186	1.53%	0.240%
25	21.454	3119	3126	3138	rVV6	65702	218182	4.84%	0.756%
26	21.701	3163	3168	3177	rBV3	186977	317515	7.04%	1.100%
27	21.795	3177	3184	3194	rVB	661130	1151143	25.52%	3.989%
28	21.954	3206	3211	3218	rVB	63705	98440	2.18%	0.341%
29	22.059	3224	3229	3238	rBV5	36487	78616	1.74%	0.272%
30	22.576	3312	3317	3324	rVB3	83473	133344	2.96%	0.462%
31	23.058	3393	3399	3405	rBV	51002	99604	2.21%	0.345%
32	23.293	3433	3439	3448	rVB	114696	215538	4.78%	0.747%
33	23.493	3466	3473	3479	rBV	49293	95543	2.12%	0.331%
34	24.128	3574	3581	3590	rVB3	104029	225603	5.00%	0.782%

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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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35	25.150	3740	3755	3770	rBV3	311911	1209676	26.82%	4.192%
36	26.296	3941	3950	3959	rBV3	57295	170953	3.79%	0.592%
37	27.712	4182	4191	4204	rBV7	31602	109887	2.44%	0.381%

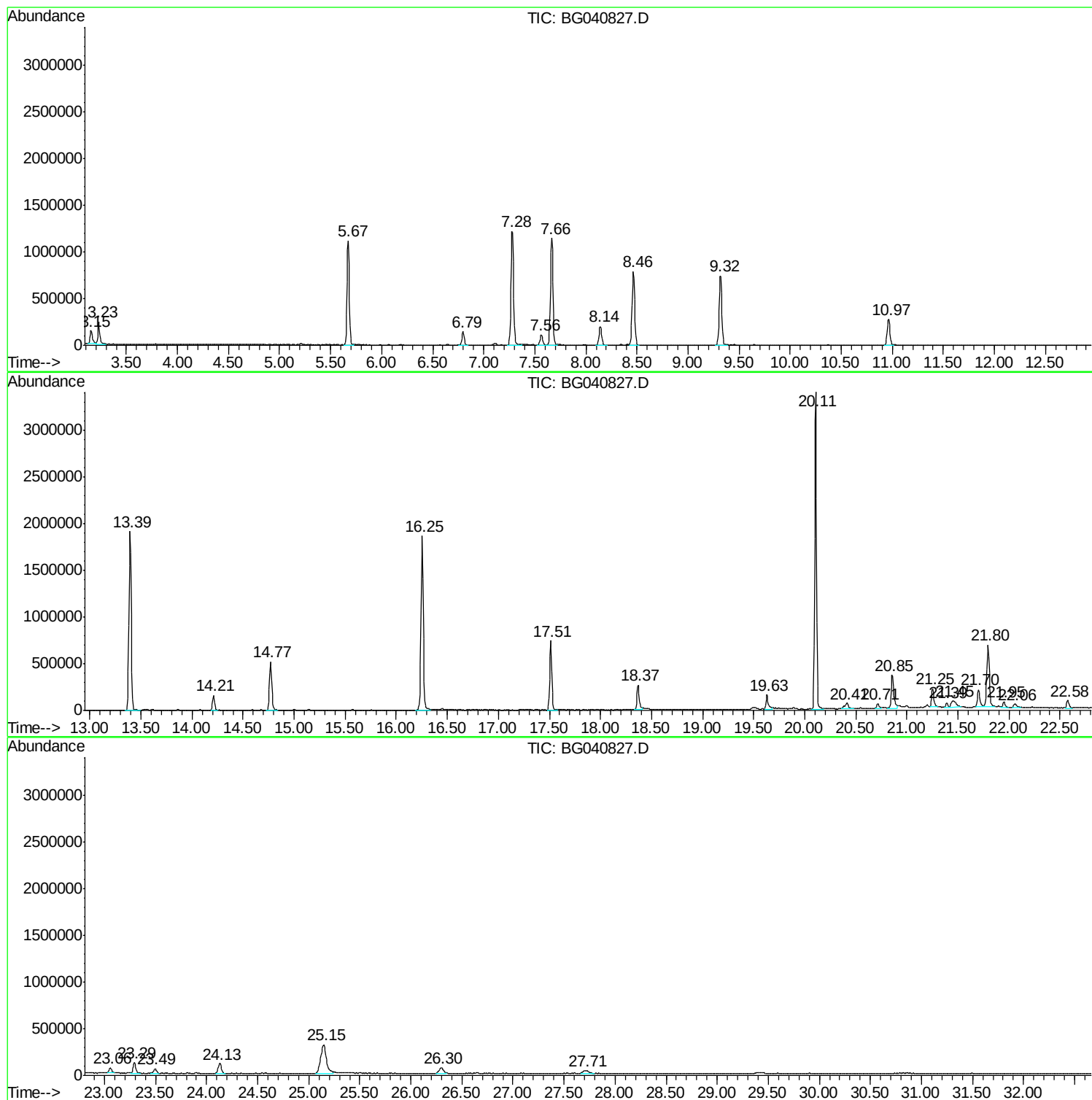
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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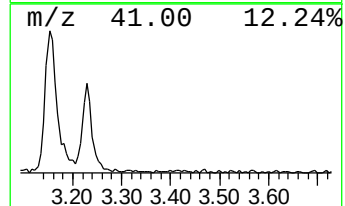
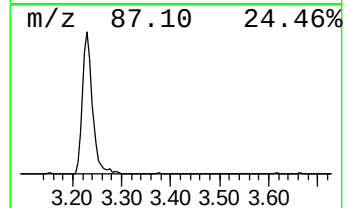
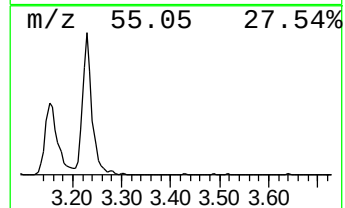
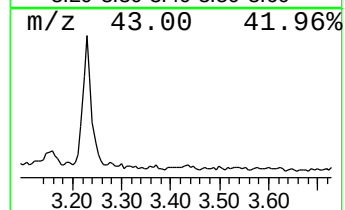
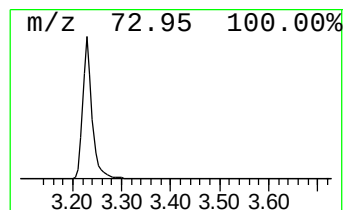
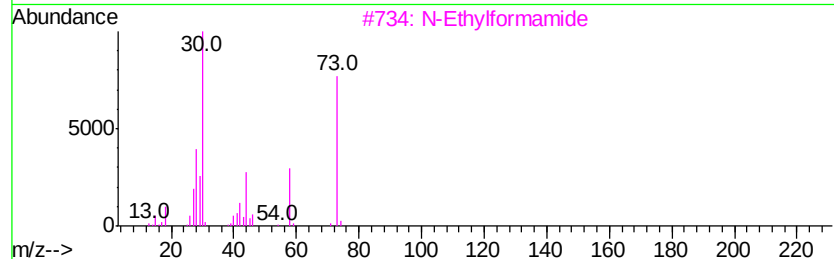
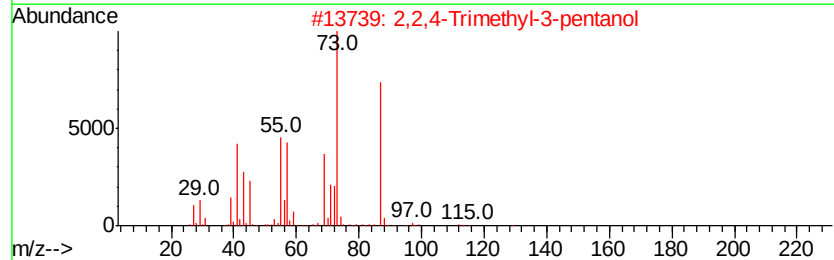
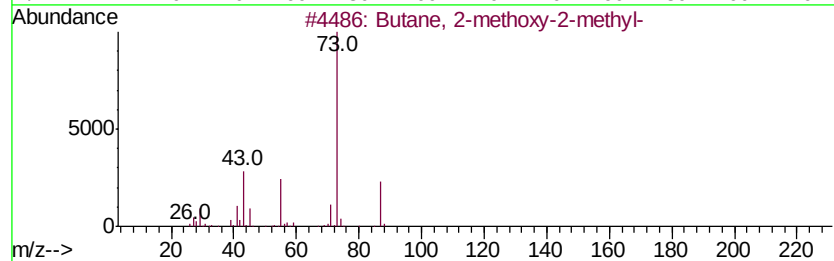
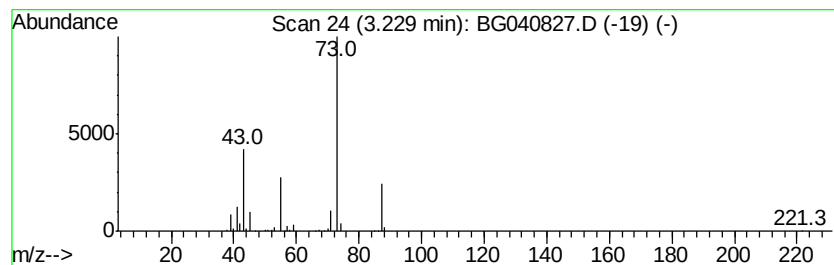
Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG042319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
3.23	18.50 ng	322532	1,4-Dichlorobenzene-d4	8.14	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	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	N-Ethylformamide	73	C3H7NO	000627-45-2	9
4	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
5	Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	9



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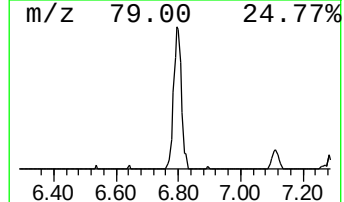
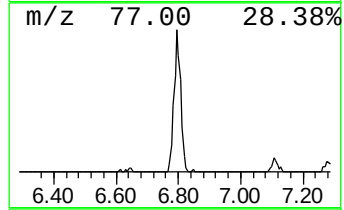
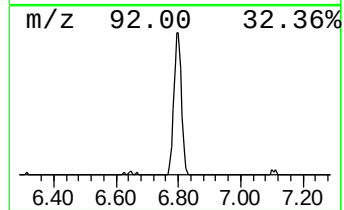
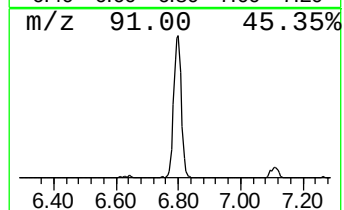
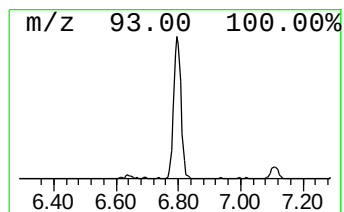
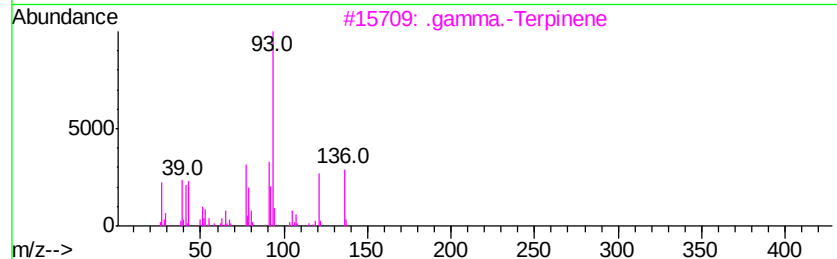
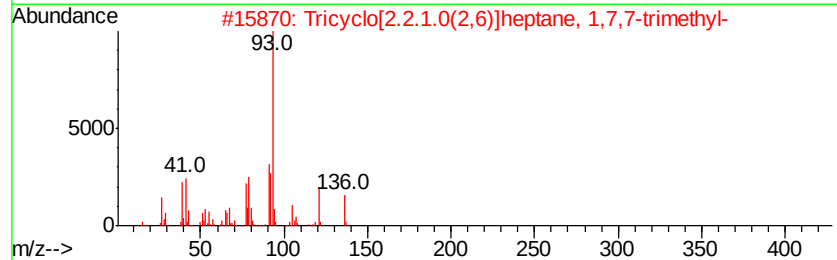
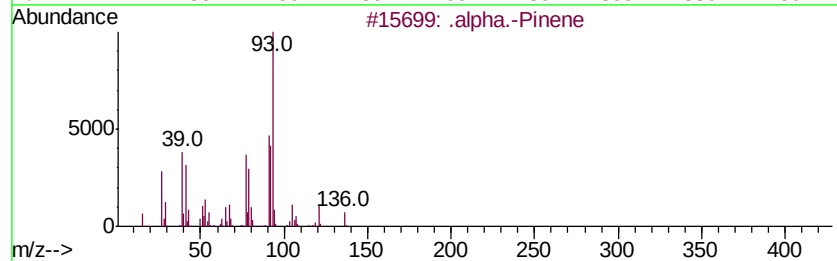
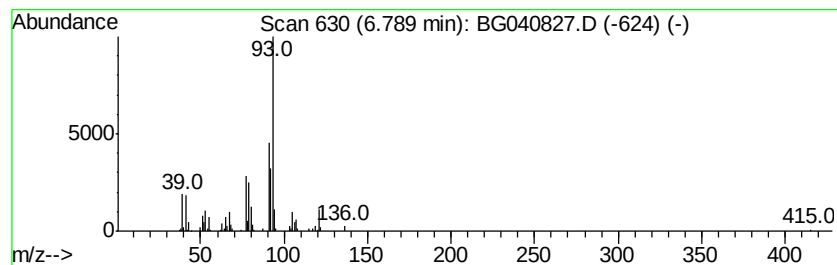
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG042319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
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Peak Number 3 .alpha.-Pinene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.79	13.54 ng	236063	1,4-Dichlorobenzene-d4	8.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.alpha.-Pinene	136	C10H16	000080-56-8	95
2		Tricyclo[2.2.1.0(2,6)]heptane, 1...	136	C10H16	000508-32-7	91
3		.gamma.-Terpinene	136	C10H16	000099-85-4	91
4		1,3,6-Octatriene, 3,7-dimethyl-,...	136	C10H16	003338-55-4	91
5		(1R)-2,6,6-Trimethylbicyclo[3.1....	136	C10H16	007785-70-8	91



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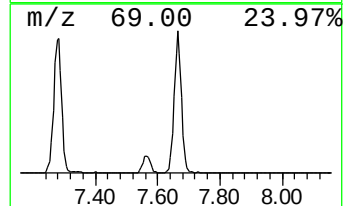
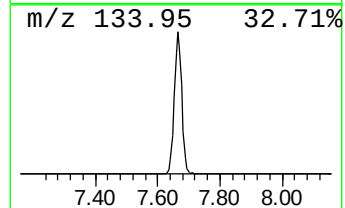
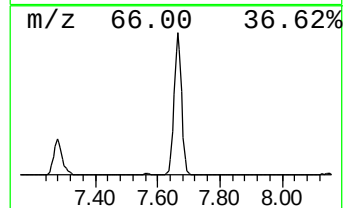
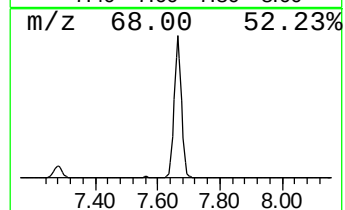
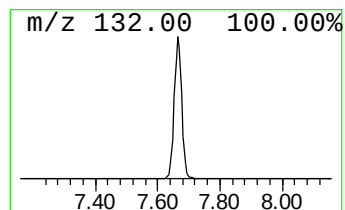
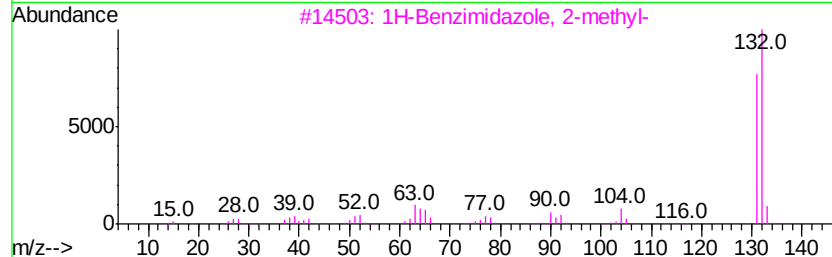
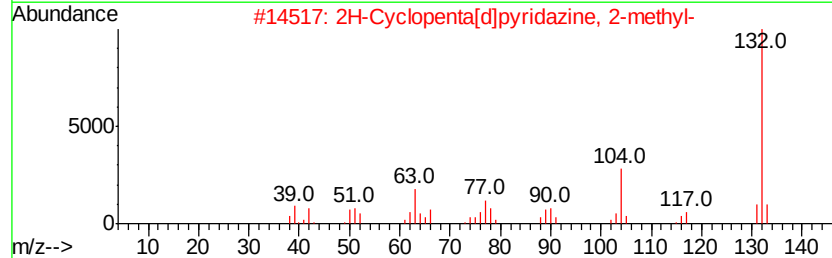
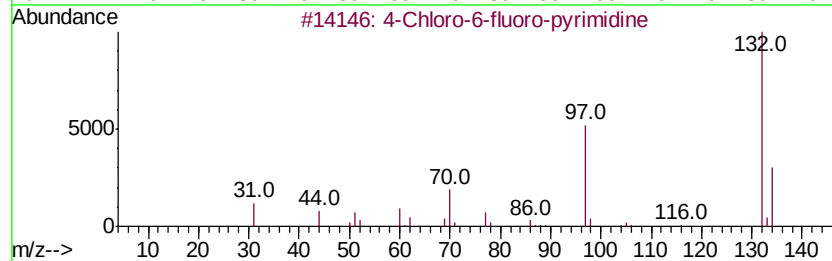
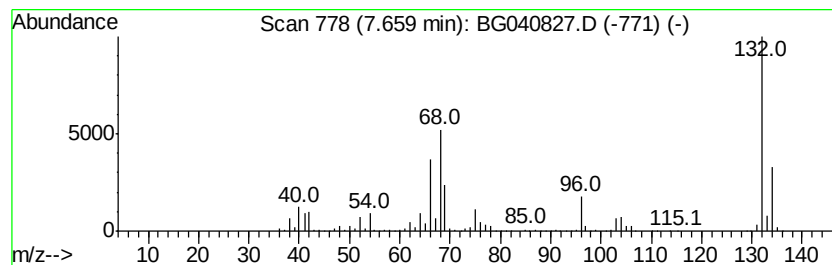
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TIC Library : C:\Database\NIST11.L
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Peak Number 4 unknown7.66 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.66	111.17 ng	1937780	1,4-Dichlorobenzene-d4	8.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	22
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	14



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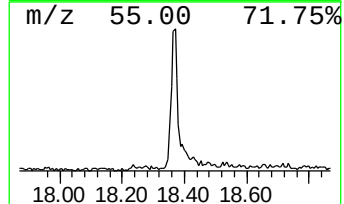
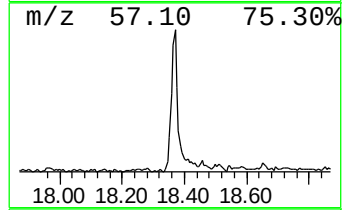
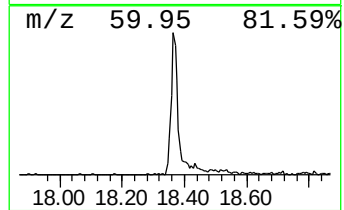
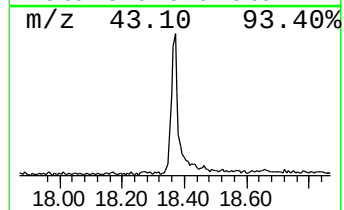
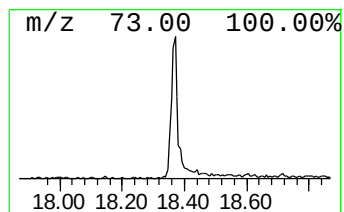
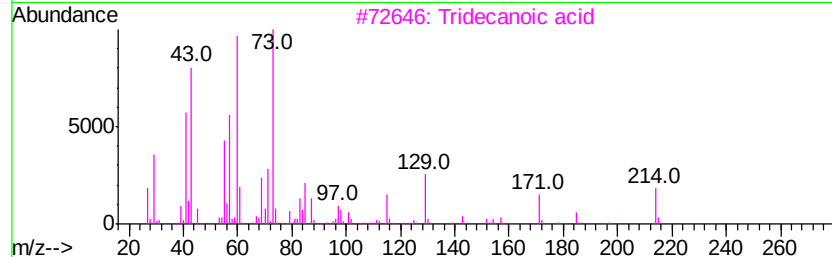
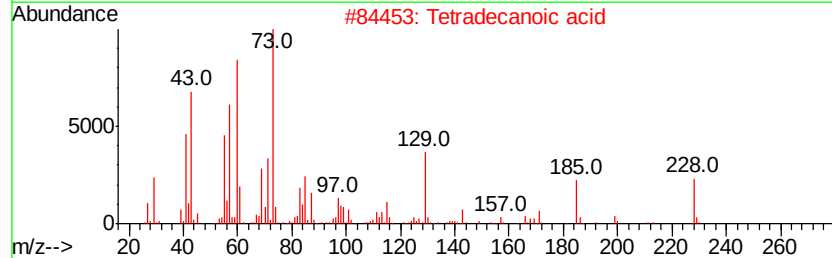
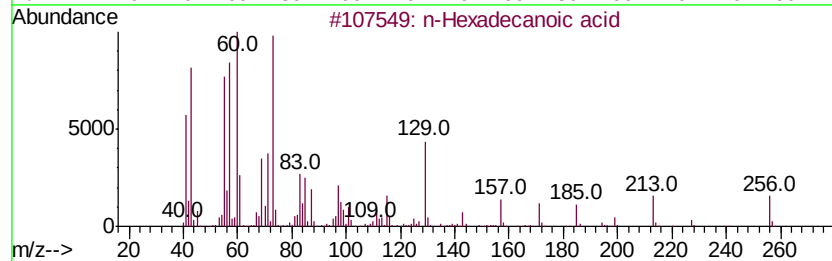
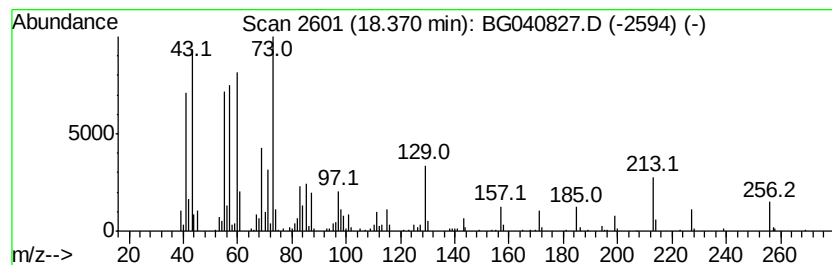
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TIC Library : C:\Database\NIST11.L
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Peak Number 6 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.37	7.33 ng	412494	Phenanthrene-d10	17.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	94
3		Tridecanoic acid	214	C13H26O2	000638-53-9	90
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	81
5		Dodecanoic acid	200	C12H24O2	000143-07-7	72



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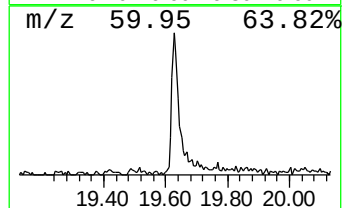
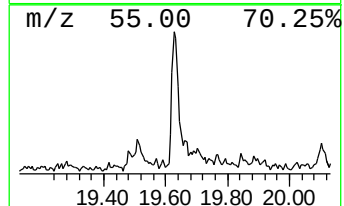
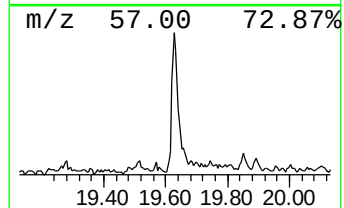
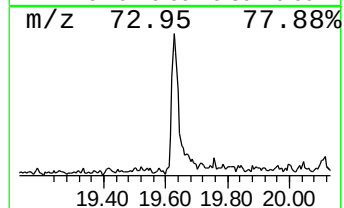
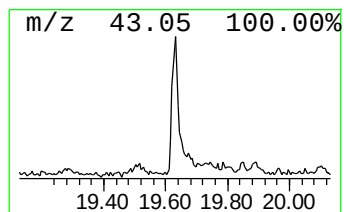
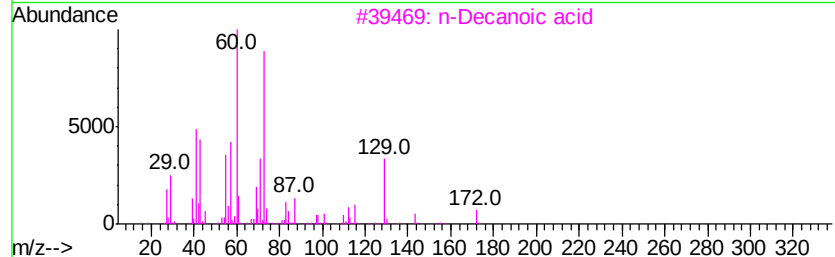
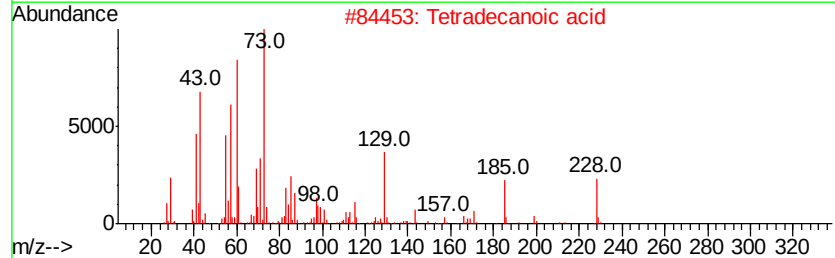
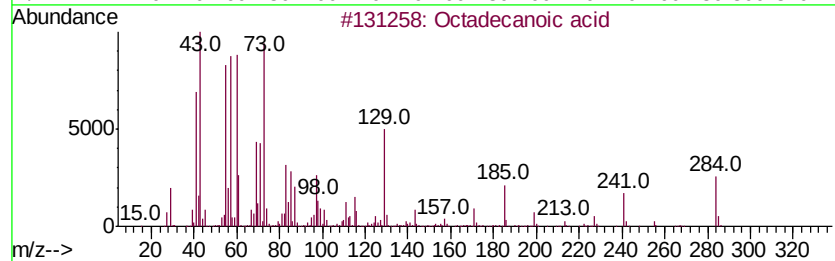
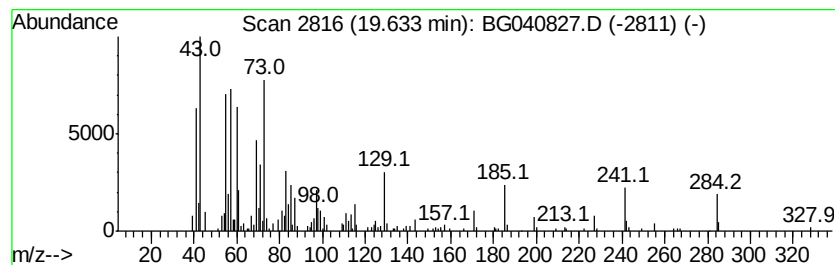
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Peak Number 7 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.63	4.64 ng	261170	Phenanthrene-d10	17.51

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	98
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	72
3			n-Decanoic acid	172	C10H20O2	000334-48-5	64
4			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	64
5			Tridecanoic acid	214	C13H26O2	000638-53-9	58



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Data File : BG040827.D
Acq On : 9 May 2019 22:18
Operator : HP/JU
Sample : K2762-16
Misc :
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Instrument :
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ClientSampleId :
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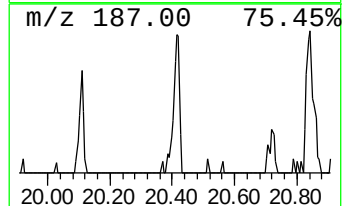
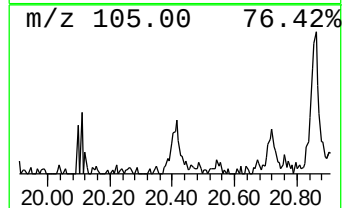
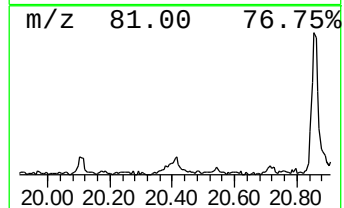
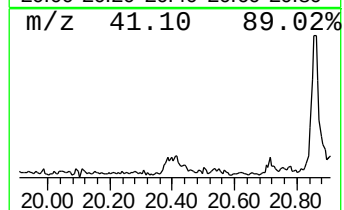
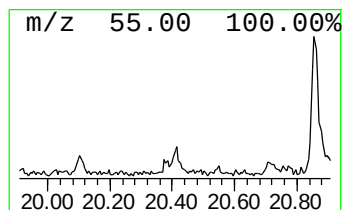
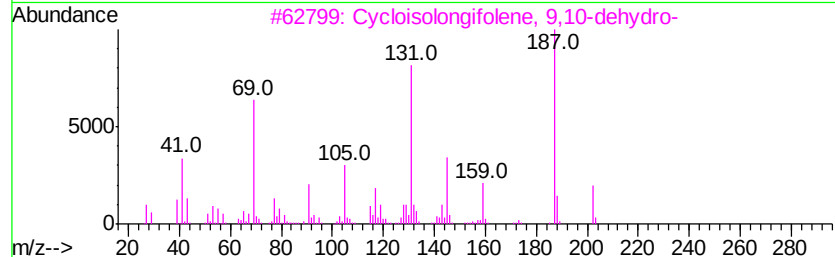
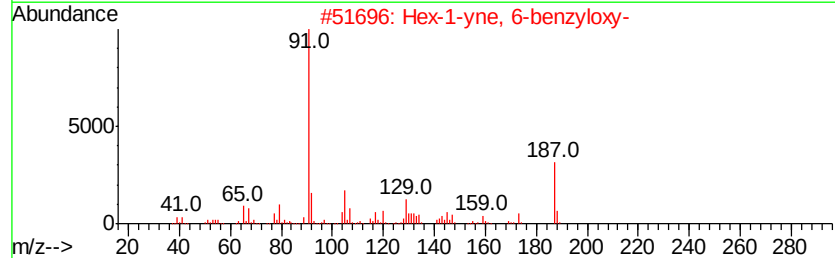
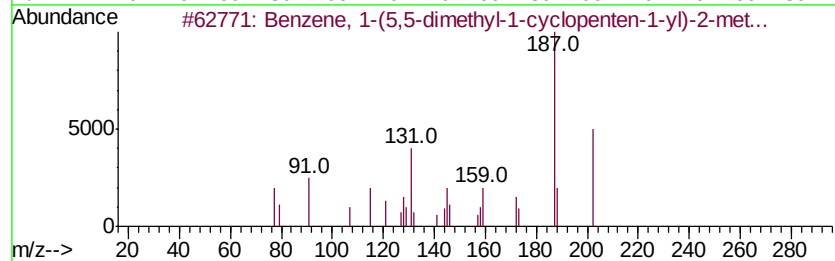
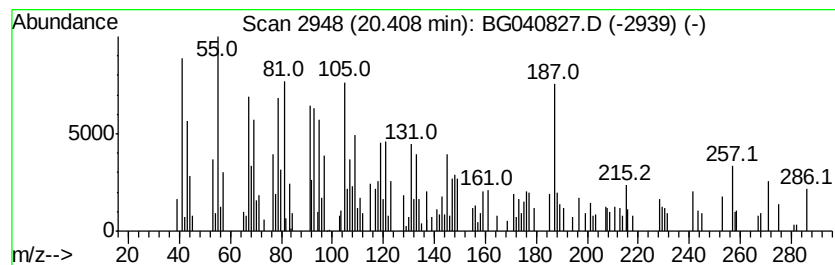
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown20.41 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.41	2.74 ng	157628	Chrysene-d12	21.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-(5,5-dimethyl-1-cyclo...	202	C14H18O	039877-93-5	46
2		Hex-1-yne, 6-benzoyloxy-	188	C13H16O	060789-55-1	25
3		Cycloisolongifolene, 9,10-dehydro-	202	C15H22	1000156-81-6	18
4		9-Borabicyclo[3.3.1]nonane, 9-(1...	164	C11H21B	051475-51-5	14
5		N-Furfuryl-N-methylaniline	187	C12H13NO	105909-65-7	11



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050919\
Data File : BG040827.D
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ClientSampleId :
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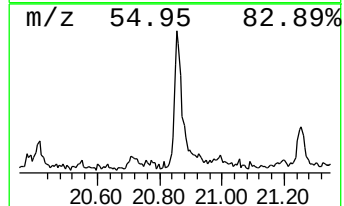
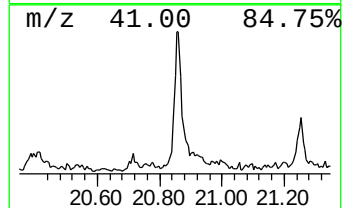
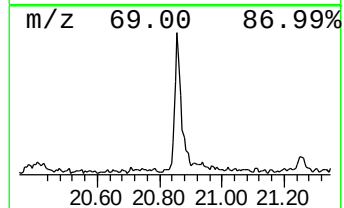
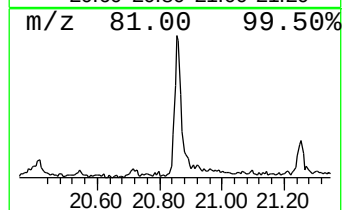
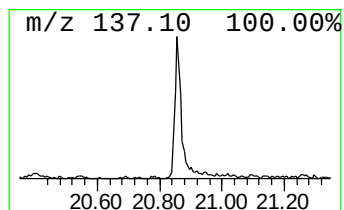
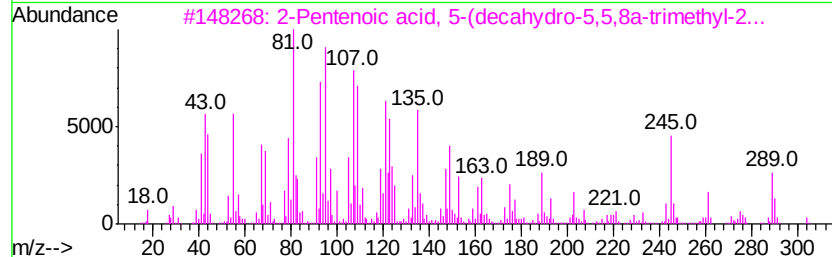
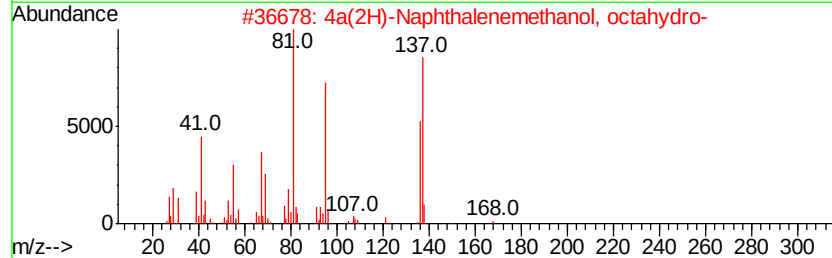
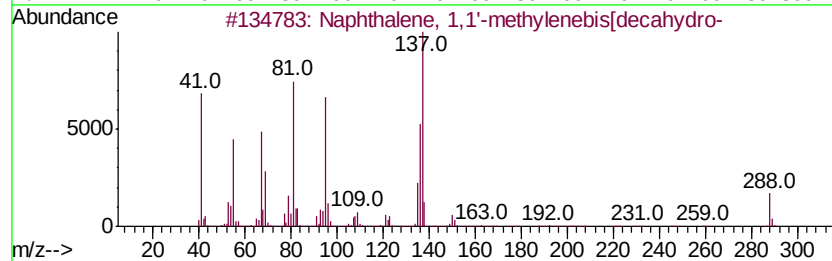
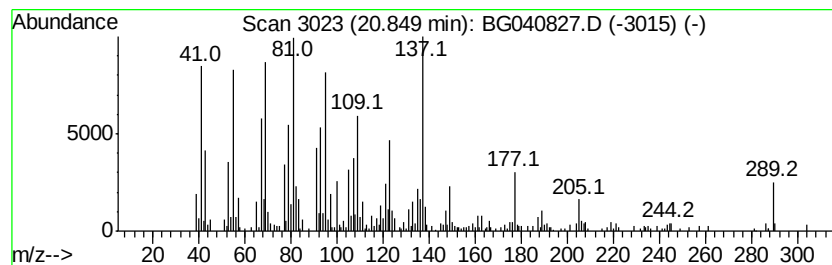
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown20.85 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.85	10.99 ng	632757	Chrysene-d12	21.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,1'-methylenebis[d...	288	C21H36	055125-02-5	46
2		4a(2H)-Naphthalenemethanol, octa...	168	C11H20O	099992-19-5	45
3		2-Pentenoic acid, 5-(decahydro-5...	304	C20H32O2	024470-48-2	42
4		cis, cis-3-Ethylbicyclo[4.4.0]de...	166	C12H22	066660-42-2	38
5		2-Cyclohexene-1-carboxaldehyde, ...	220	C15H24O	056772-07-7	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050919\
Data File : BG040827.D
Acq On : 9 May 2019 22:18
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Sample : K2762-16
Misc :
ALS Vial : 8 Sample Multiplier: 1

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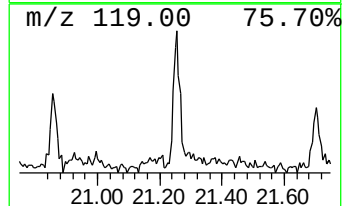
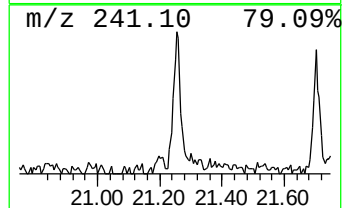
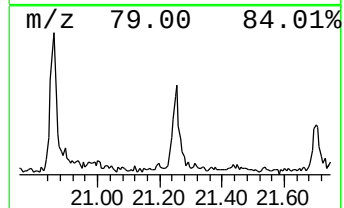
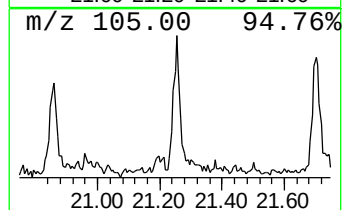
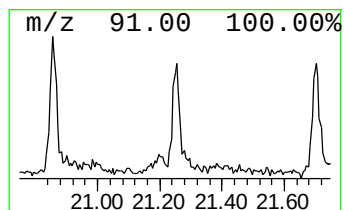
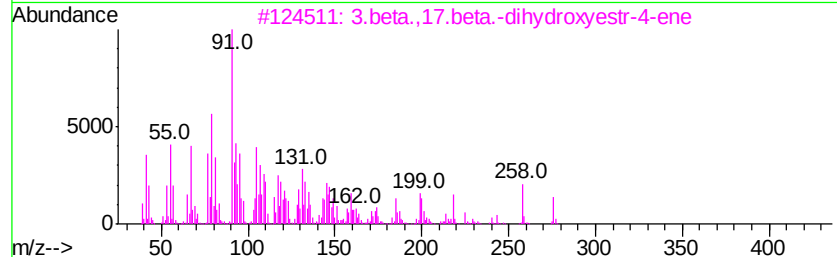
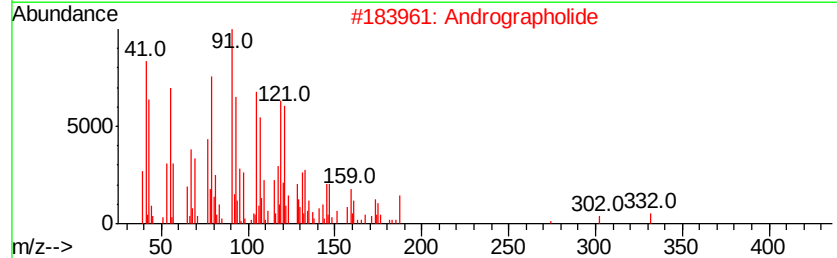
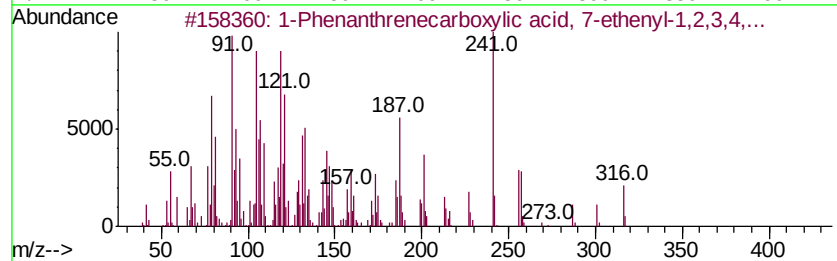
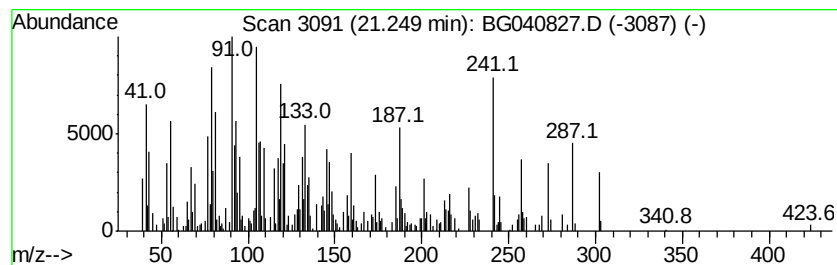
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown21.25 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.25	5.17 ng	297633	Chrysene-d12	21.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Phenanthrenecarboxylic acid, 7...	316	C21H32O2	001686-62-0	35
2		Andrographolide	350	C20H30O5	005508-58-7	32
3		3.beta.,17.beta.-dihydroxvestr-4...	276	C18H28O2	019793-20-5	32
4		3.alpha.,17.beta.-dihydroxvestr-...	276	C18H28O2	035950-87-9	12
5		3-Phenylpropionic acid, 2,7-dime...	284	C19H24O2	1000299-17-7	12



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050919\
Data File : BG040827.D
Acq On : 9 May 2019 22:18
Operator : HP/JU
Sample : K2762-16
Misc :
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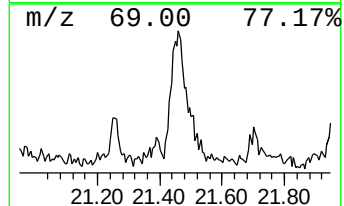
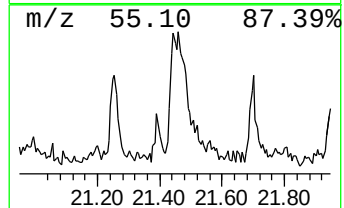
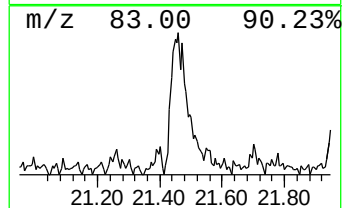
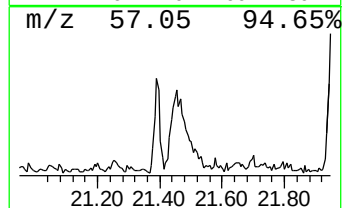
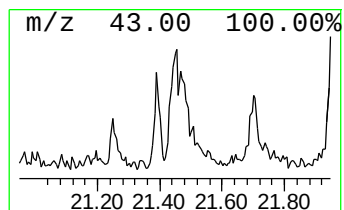
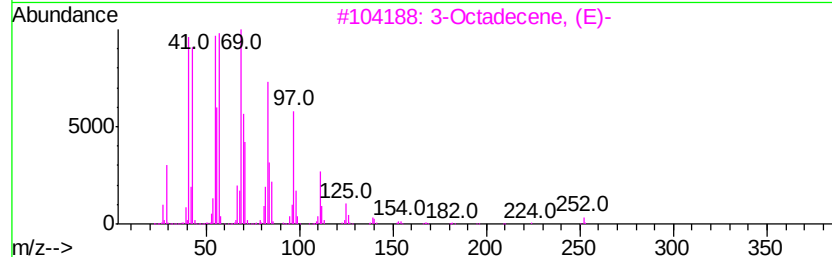
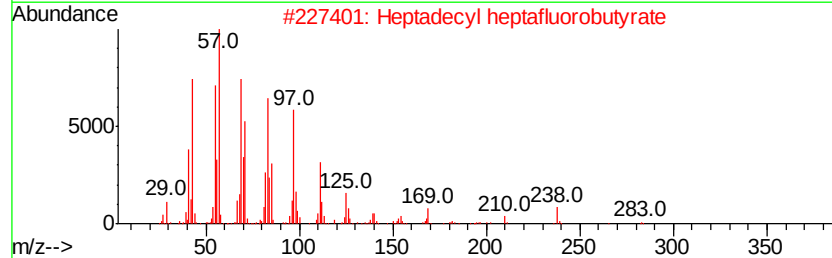
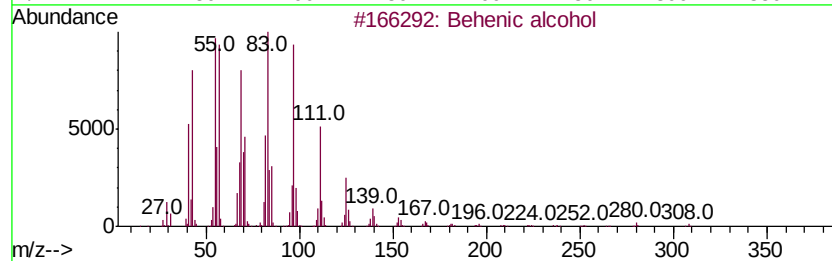
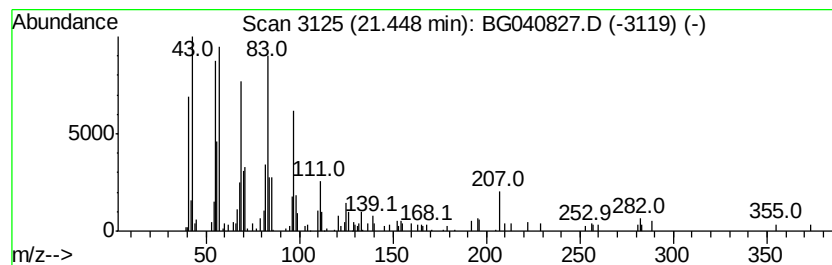
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Behenic alcohol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.45	3.79 ng	218182	Chrysene-d12	21.80	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Behenic alcohol	326	C22H46O	000661-19-8	91
2	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	90
3	3-Octadecene, (E)-	252	C18H36	007206-19-1	90
4	Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	87
5	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050919\
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Sample : K2762-16
Misc :
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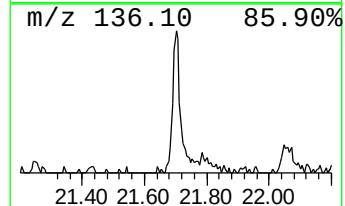
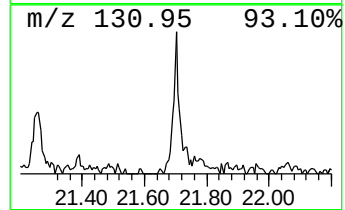
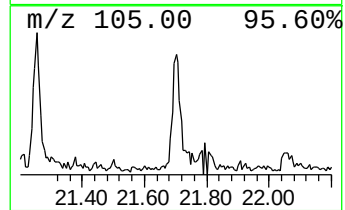
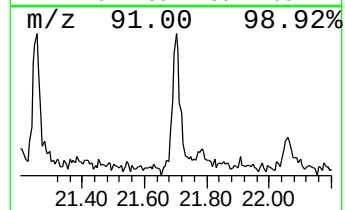
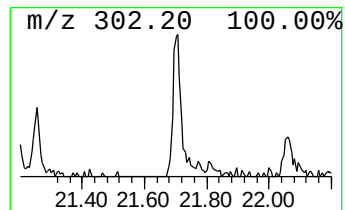
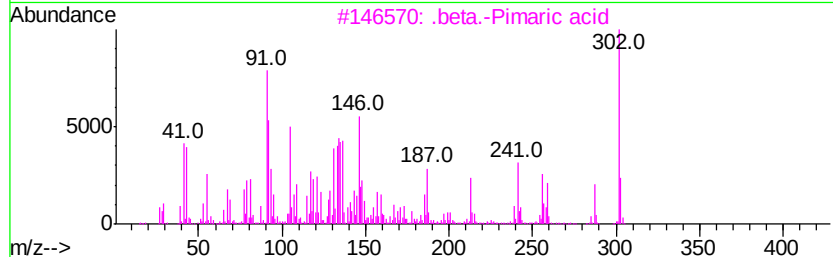
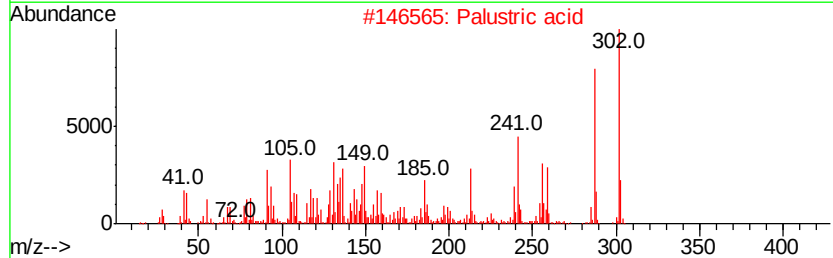
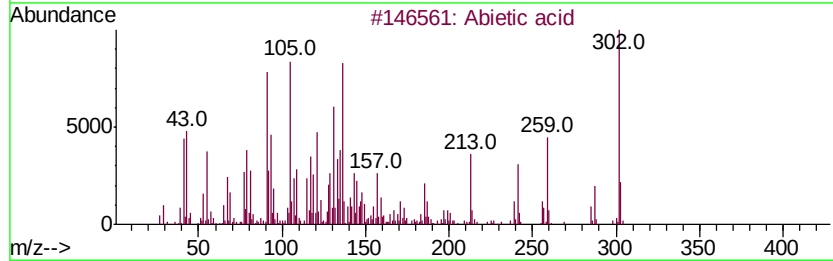
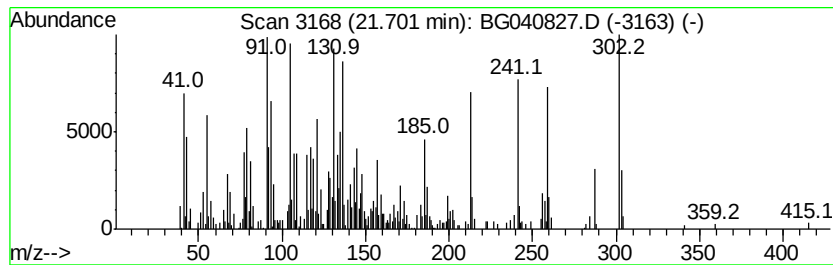
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Abietic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
21.70	5.52 ng	317515	Chrysene-d12	21.80		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Abietic acid		302	C20H30O2	000514-10-3	98
2	Palustric acid		302	C20H30O2	001945-53-5	64
3	.beta.-Pimaric acid		302	C20H30O2	000079-54-9	27
4	Methyl abietate		316	C21H32O2	000127-25-3	22
5	4-(1-Naphthyl)-5-phenyl-2-thiazo...		302	C19H14N2S	109693-78-9	15



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050919\
Data File : BG040827.D
Acq On : 9 May 2019 22:18
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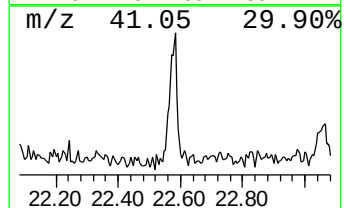
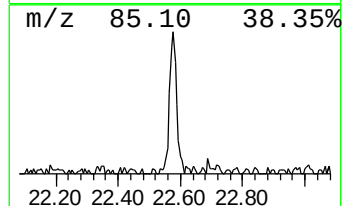
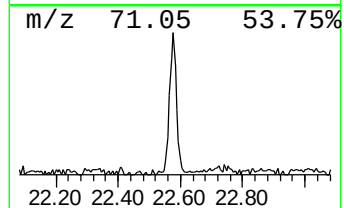
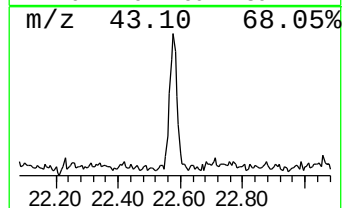
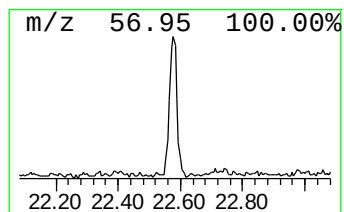
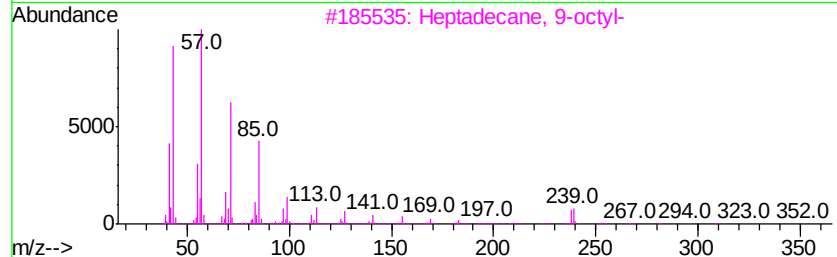
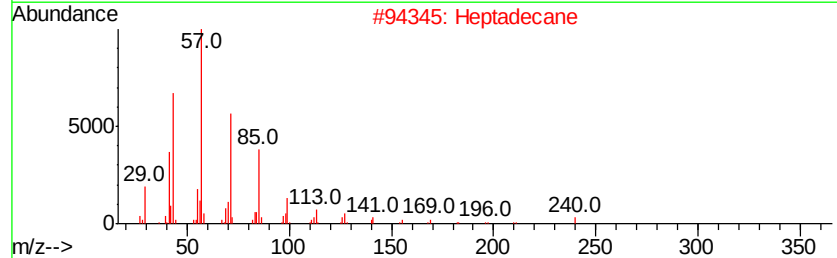
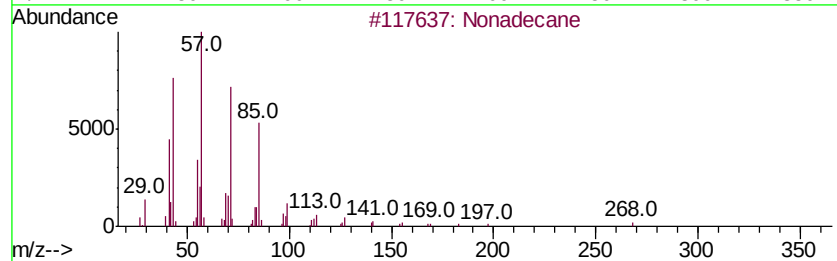
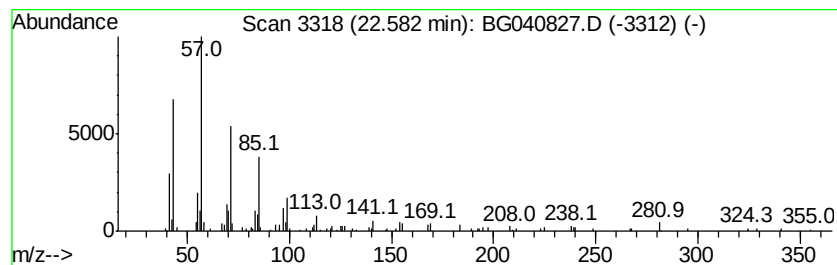
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Nonadecane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.58	2.32 ng	133344	Chrysene-d12	21.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	98
2		Heptadecane	240	C17H36	000629-78-7	96
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
4		Octadecane	254	C18H38	000593-45-3	90
5		Hexadecane	226	C16H34	000544-76-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050919\
Data File : BG040827.D
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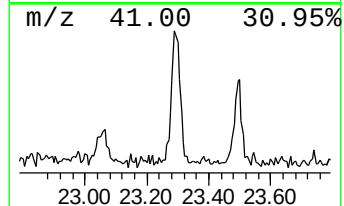
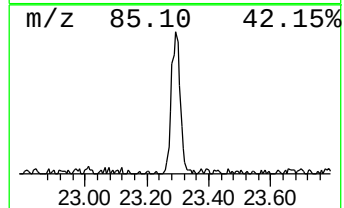
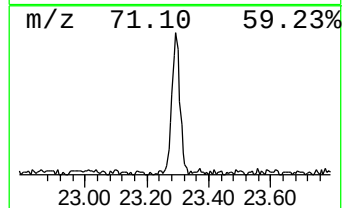
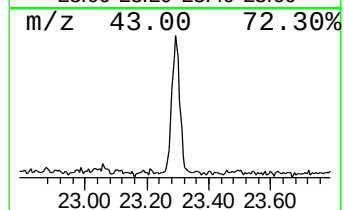
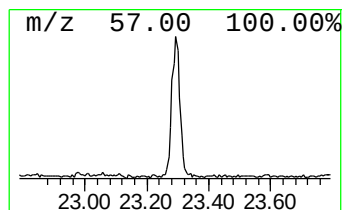
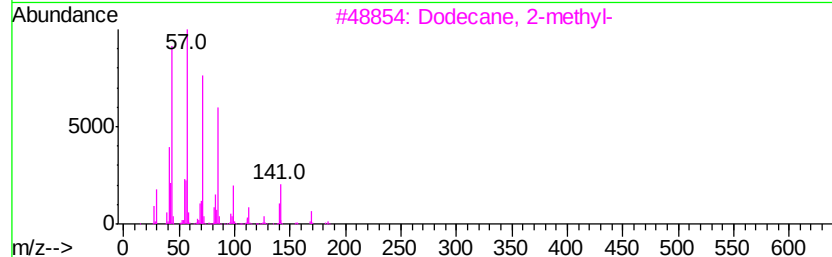
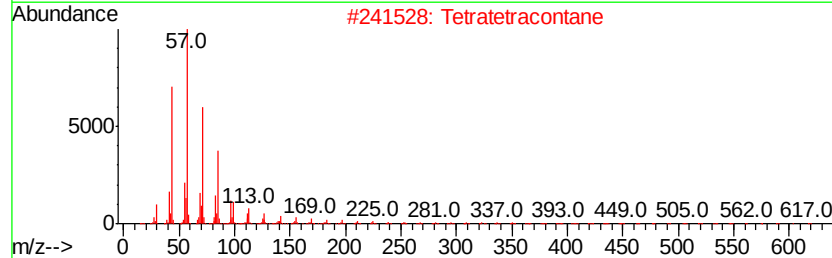
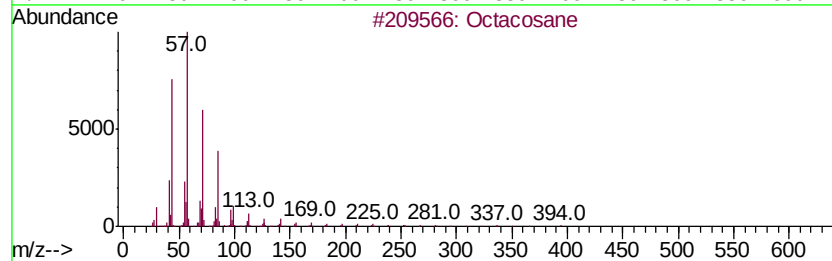
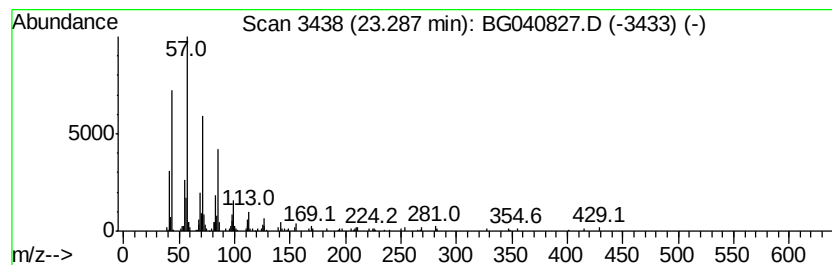
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Octacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.29	3.74 ng	215538	Chrysene-d12	21.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Tetratetracontane	619	C44H90	007098-22-8	91
3		Dodecane, 2-methyl-	184	C13H28	001560-97-0	90
4		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	87
5		Hexatriacontane	507	C36H74	000630-06-8	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050919\
Data File : BG040827.D
Acq On : 9 May 2019 22:18
Operator : HP/JU
Sample : K2762-16
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
P021-SS003-0612-01

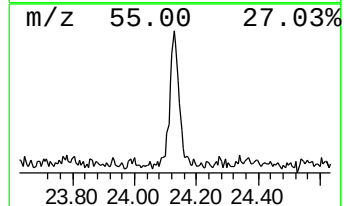
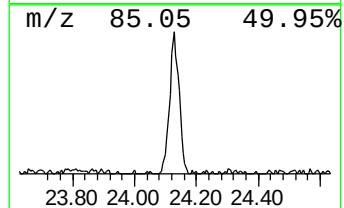
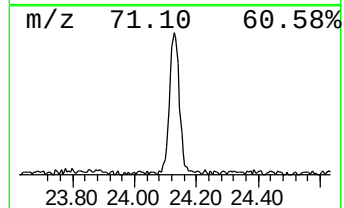
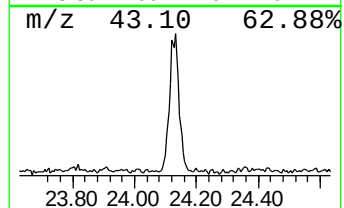
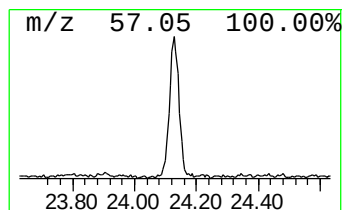
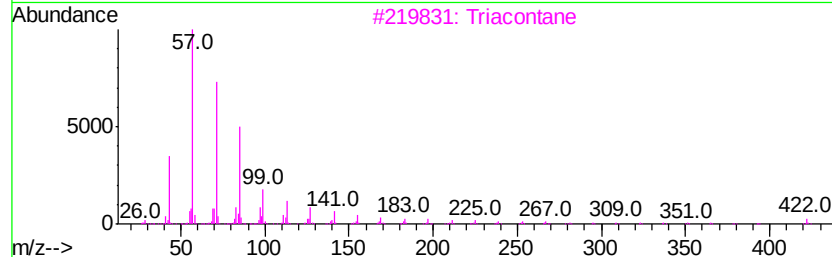
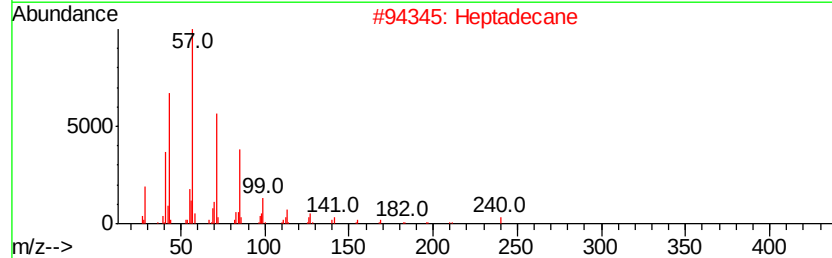
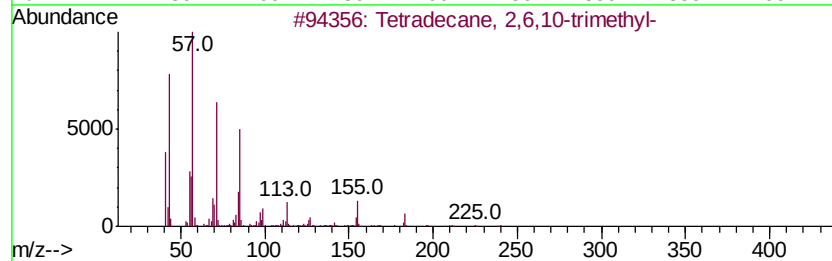
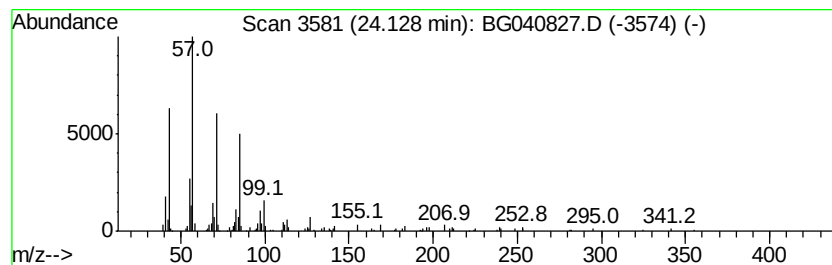
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Tetradecane, 2,6,10-trimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.13	3.73 ng	225603	Perylene-d12	25.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	90
2			Heptadecane	240	C17H36	000629-78-7	87
3			Triacontane	422	C30H62	000638-68-6	87
4			5-Ethyl-5-methylnonadecane	310	C22H46	1000360-42-7	80
5			Tetratetracontane	619	C44H90	007098-22-8	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050919\
Data File : BG040827.D
Acq On : 9 May 2019 22:18
Operator : HP/JU
Sample : K2762-16
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
P021-SS003-0612-01

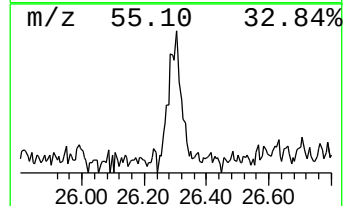
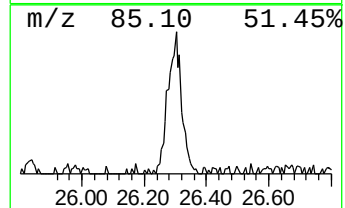
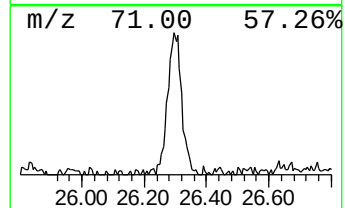
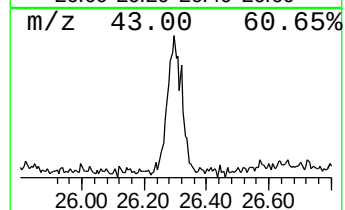
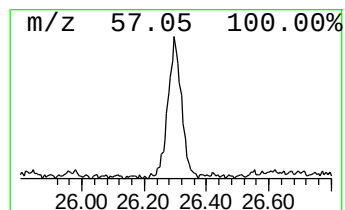
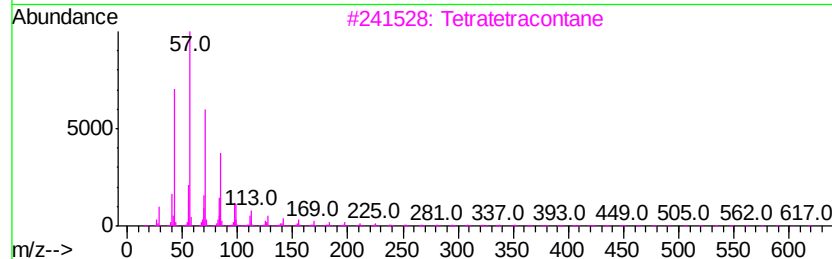
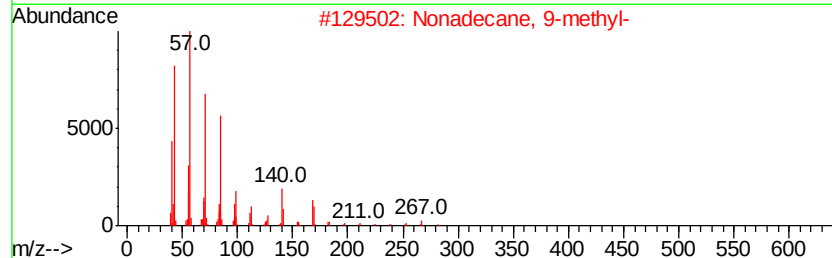
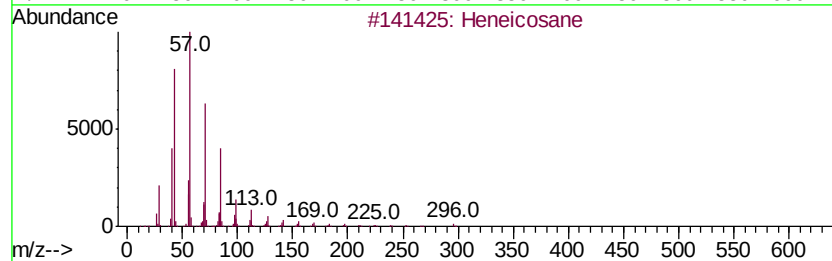
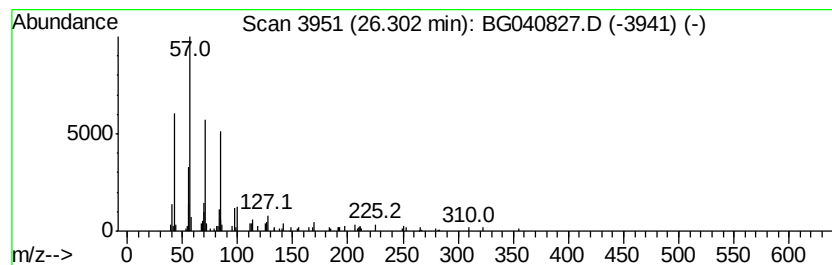
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Heneicosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.30	2.83 ng	170953	Perylene-d12	25.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	96
2			Nonadecane, 9-methyl-	282	C20H42	013287-24-6	87
3			Tetratetracontane	619	C44H90	007098-22-8	87
4			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
5			Docosane, 11-butyl-	366	C26H54	013475-76-8	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG050919\
 Data File : BG040827.D
 Acq On : 9 May 2019 22:18
 Operator : HP/JU
 Sample : K2762-16
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 P021-SS003-0612-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG042319.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...	3.23	18.5	ng	322532	1	8.14	348628	20.0
.alpha.-Pinene	6.79	13.5	ng	236063	1	8.14	348628	20.0
unknown7.66	7.66	111.2	ng	1937780	1	8.14	348628	20.0
n-Hexadecanoic acid	18.37	7.3	ng	412494	4	17.51	1125730	20.0
Octadecanoic acid	19.63	4.6	ng	261170	4	17.51	1125730	20.0
unknown20.41	20.41	2.7	ng	157628	5	21.80	1151140	20.0
unknown20.85	20.85	11.0	ng	632757	5	21.80	1151140	20.0
unknown21.25	21.25	5.2	ng	297633	5	21.80	1151140	20.0
Behenic alcohol	21.45	3.8	ng	218182	5	21.80	1151140	20.0
Abietic acid	21.70	5.5	ng	317515	5	21.80	1151140	20.0
Nonadecane	22.58	2.3	ng	133344	5	21.80	1151140	20.0
Octacosane	23.29	3.7	ng	215538	5	21.80	1151140	20.0
Tetradecane, 2,6,...	24.13	3.7	ng	225603	6	25.15	1209680	20.0
Heneicosane	26.30	2.8	ng	170953	6	25.15	1209680	20.0