

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011720\
 Data File : BG044085.D
 Acq On : 17 Jan 2020 21:55
 Operator : JU
 Sample : L1162-09
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MH-719

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	5.038	326	332	346	rVB	108917	178818	3.25%	0.570%
2	5.508	406	412	433	rBV	714023	1217510	22.16%	3.883%
3	7.100	676	683	699	rBV	487145	948184	17.26%	3.024%
4	7.464	737	745	756	rBV	1367028	2468090	44.92%	7.872%
5	7.934	817	825	832	rBV	273793	487007	8.86%	1.553%
6	8.258	872	880	887	rBV	1081908	1945203	35.40%	6.204%
7	9.033	1004	1012	1015	rBV	33664	59525	1.08%	0.190%
8	9.086	1015	1021	1034	rVB	886238	1639973	29.85%	5.231%
9	10.731	1293	1301	1310	rBV	398750	739495	13.46%	2.359%
10	13.193	1712	1720	1730	rBV	2526910	4123341	75.05%	13.151%
11	14.021	1855	1861	1866	rBV	140080	216924	3.95%	0.692%
12	14.568	1947	1954	1961	rBV2	694967	1136938	20.69%	3.626%
13	14.632	1961	1965	1972	rVB	38086	61450	1.12%	0.196%
14	16.054	2200	2207	2219	rBV	2061132	3314378	60.32%	10.571%
15	17.312	2414	2421	2432	rBV	801743	1247863	22.71%	3.980%
16	18.216	2571	2575	2583	rBV4	32981	56329	1.03%	0.180%
17	19.926	2860	2866	2879	rVB	3952174	5494404	100.00%	17.524%
18	20.226	2912	2917	2926	rBV2	72934	153204	2.79%	0.489%
19	20.737	3001	3004	3008	rVB	71690	77864	1.42%	0.248%
20	21.231	3083	3088	3101	rBV	350463	620149	11.29%	1.978%
21	21.466	3124	3128	3132	rBV	50304	68327	1.24%	0.218%
22	21.560	3137	3144	3154	rBV2	754509	1296239	23.59%	4.134%
23	21.771	3175	3180	3188	rBV	146265	211961	3.86%	0.676%
24	22.365	3275	3281	3294	rBV	169688	304989	5.55%	0.973%
25	23.034	3389	3395	3406	rVB	165227	328209	5.97%	1.047%
26	23.216	3420	3426	3435	rVB	524603	977334	17.79%	3.117%
27	23.816	3522	3528	3535	rBV2	132755	271221	4.94%	0.865%
28	24.668	3664	3673	3680	rBV	464166	1269233	23.10%	4.048%
29	24.738	3681	3685	3692	rVB3	107649	232581	4.23%	0.742%
30	25.825	3863	3870	3880	rVB3	73130	206124	3.75%	0.657%

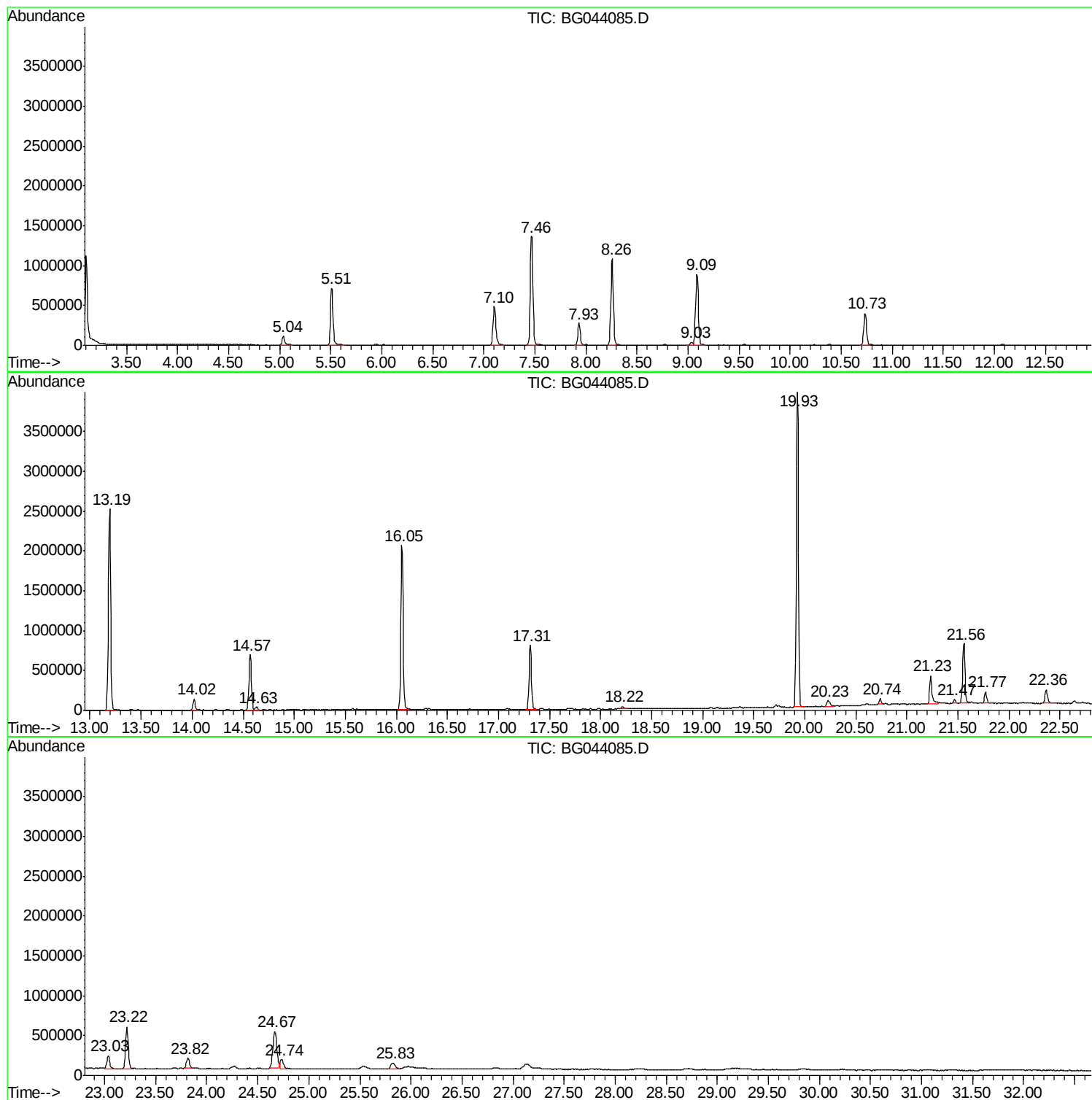
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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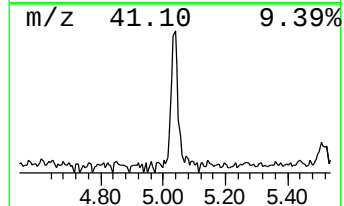
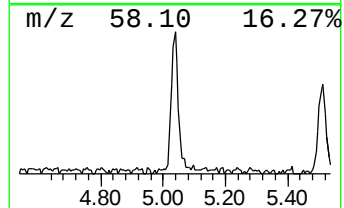
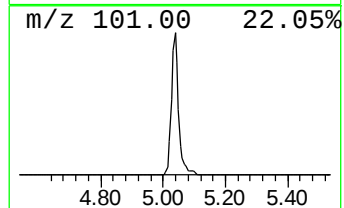
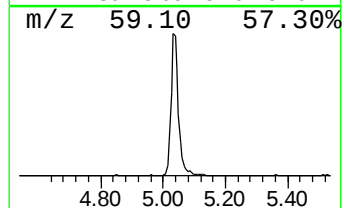
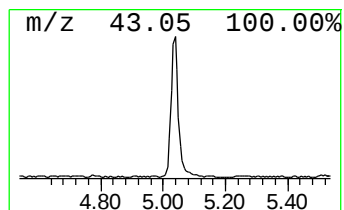
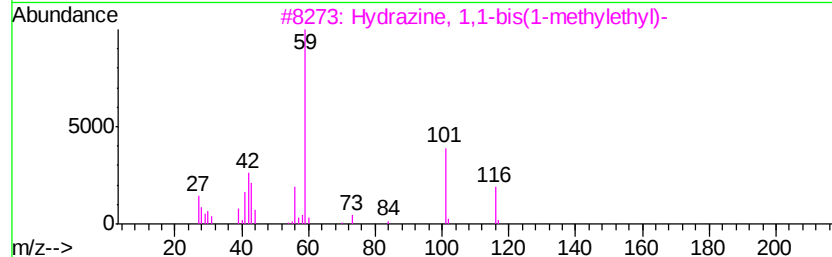
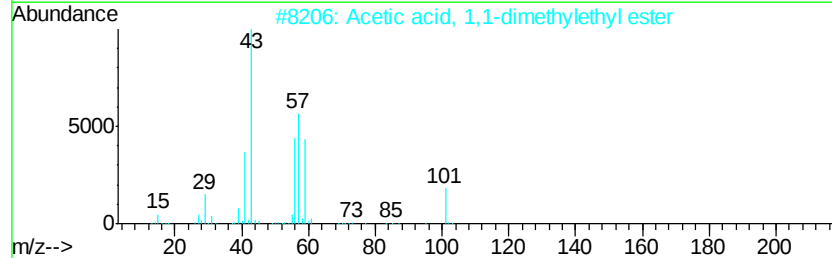
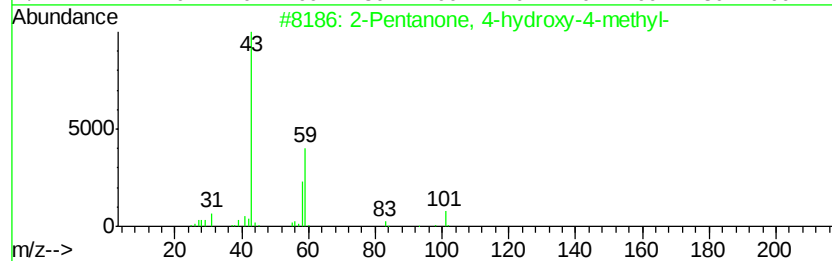
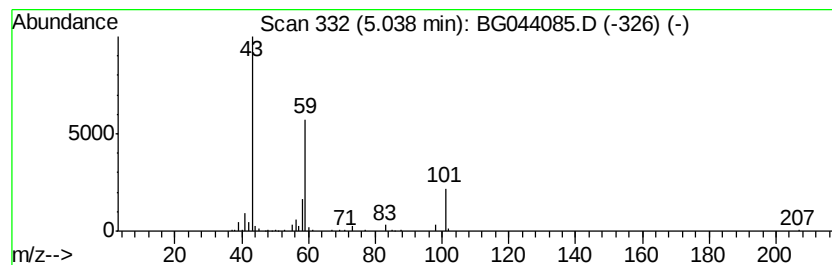
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.04	7.34 ng	178818	1,4-Dichlorobenzene-d4	7.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	33
4		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	32
5		Tetraglyme	222	C10H22O5	000143-24-8	25



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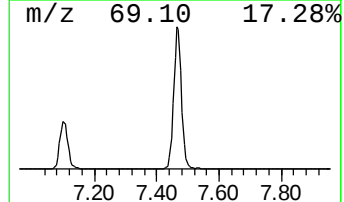
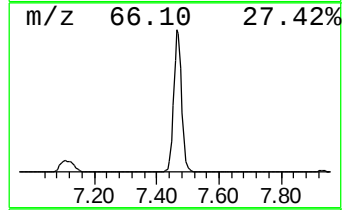
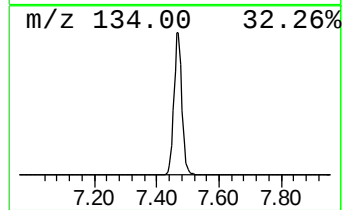
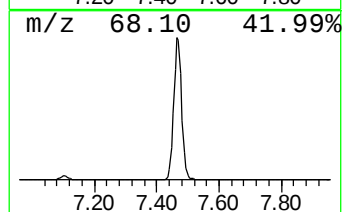
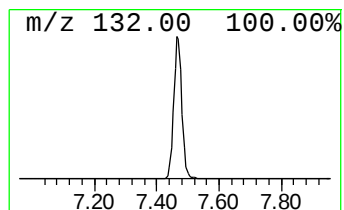
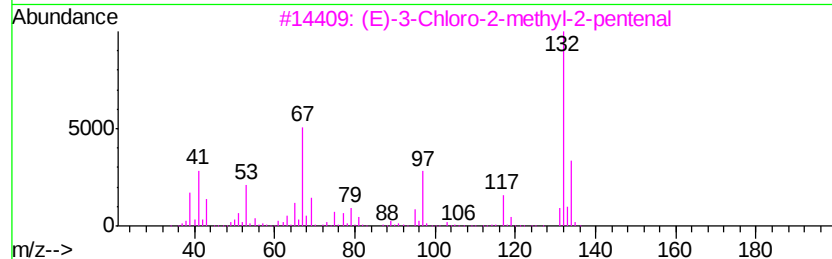
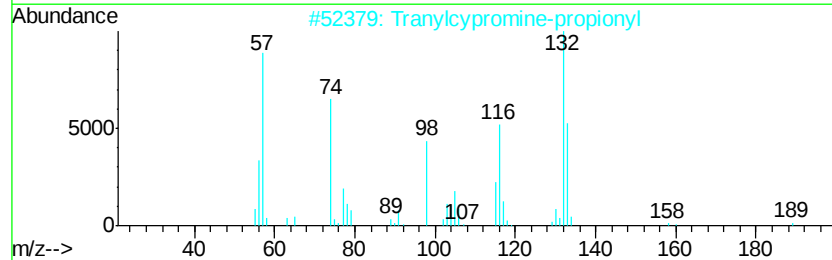
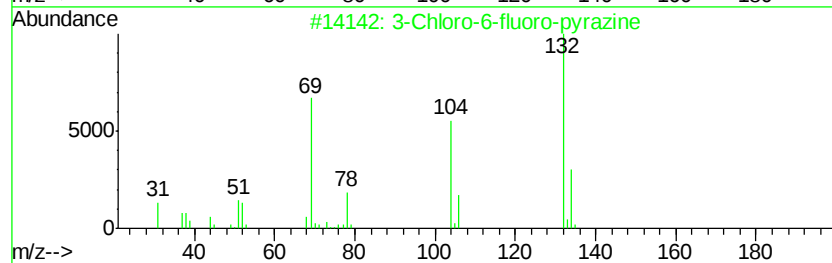
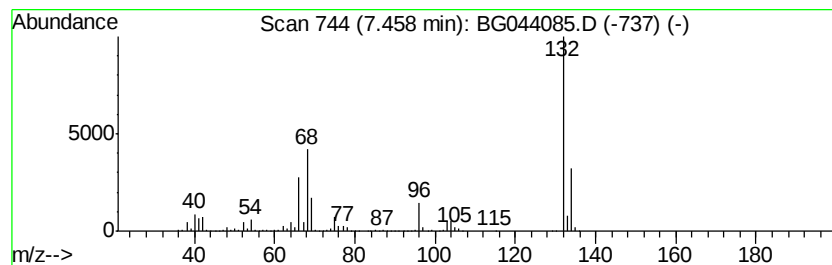
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Peak Number 2 unknown7.46 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.46	101.36 ng	2468090	1,4-Dichlorobenzene-d4	7.93

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
4			Amphetaminil	250	C17H18N2	017590-01-1	12
5			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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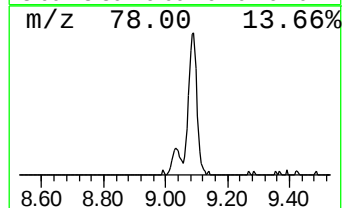
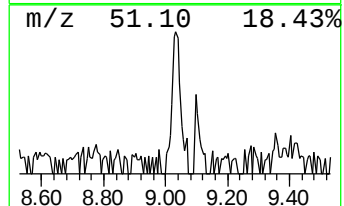
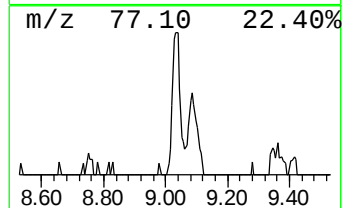
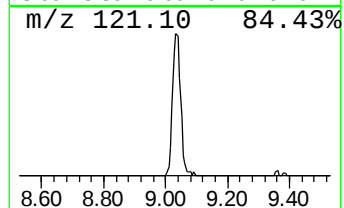
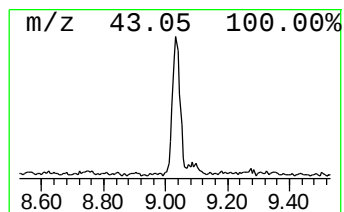
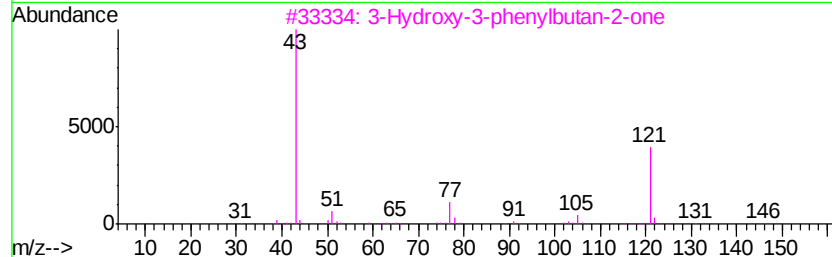
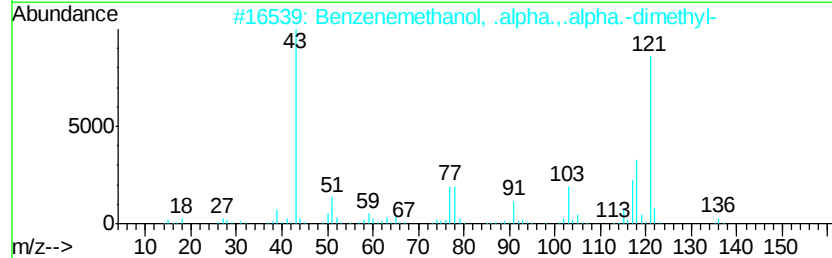
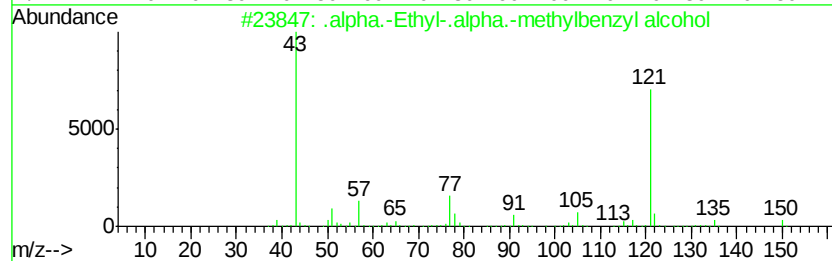
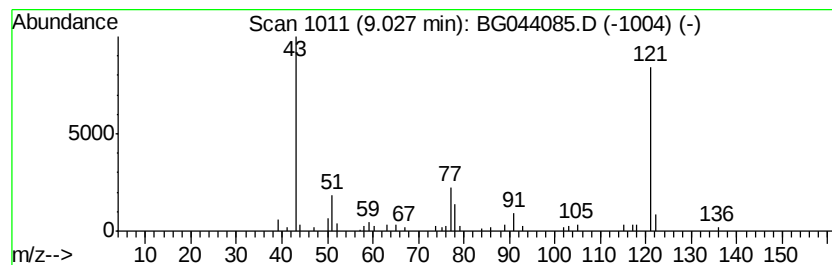
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Peak Number 4 .alpha.-Ethyl-.alpha.-methy... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.03	2.44 ng	59525	1,4-Dichlorobenzene-d4	7.93

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		.alpha.-Ethyl-.alpha.-methylbenz...	150	C10H14O	001565-75-9	78
2		Benzenemethanol, .alpha.,.alpha....	136	C9H12O	000617-94-7	72
3		3-Hydroxy-3-phenylbutan-2-one	164	C10H12O2	1000368-45-5	64
4		Benzenamine, 2,4-dimethyl-	121	C8H11N	000095-68-1	58
5		Atrolactic acid	166	C9H10O3	000515-30-0	56



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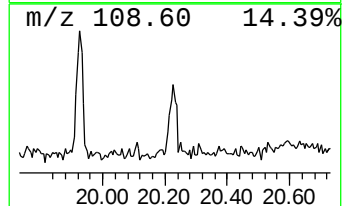
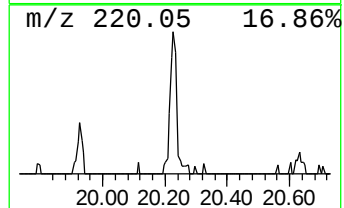
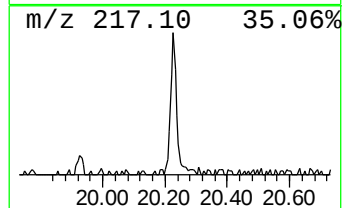
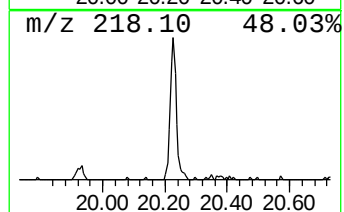
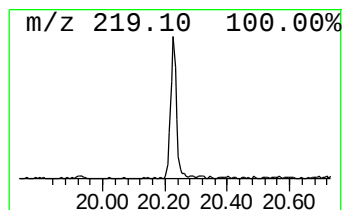
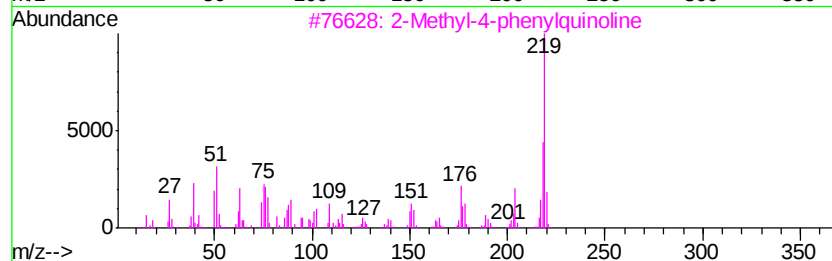
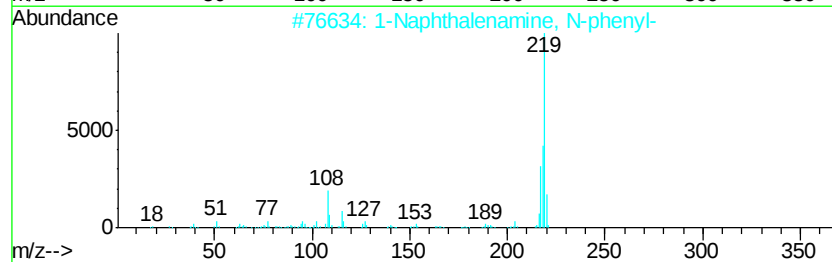
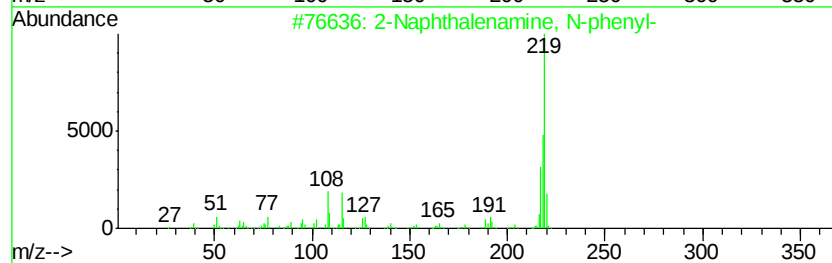
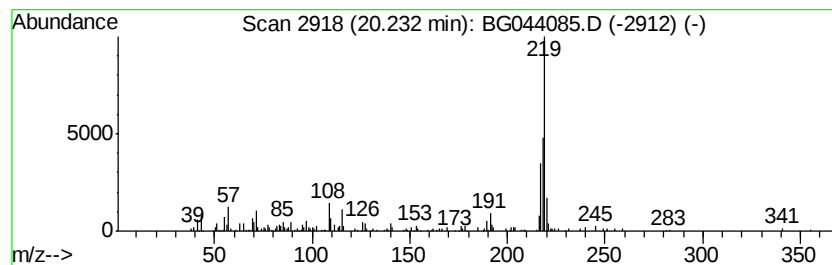
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Peak Number 5 2-Naphthalenamine, N-phenyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.23	2.36 ng	153204	Chrysene-d12	21.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Naphthalenamine, N-phenyl-	219	C16H13N	000135-88-6	92
2		1-Naphthalenamine, N-phenyl-	219	C16H13N	000090-30-2	70
3		2-Methyl-4-phenylquinoline	219	C16H13N	001721-92-2	64
4		Quinoline, 4-methyl-2-phenyl-	219	C16H13N	004789-76-8	59
5		1H-Pyrrole, 2,4-diphenyl-	219	C16H13N	003274-56-4	55



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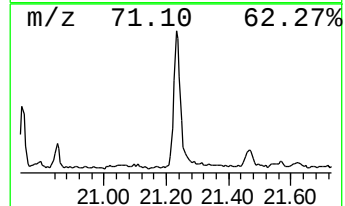
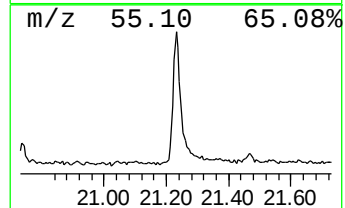
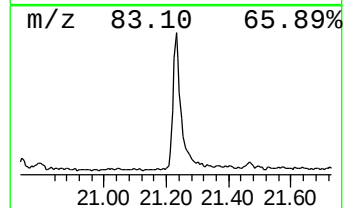
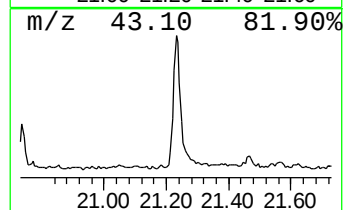
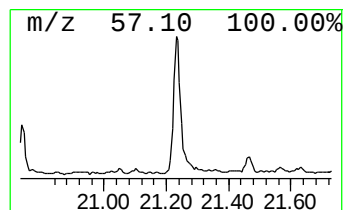
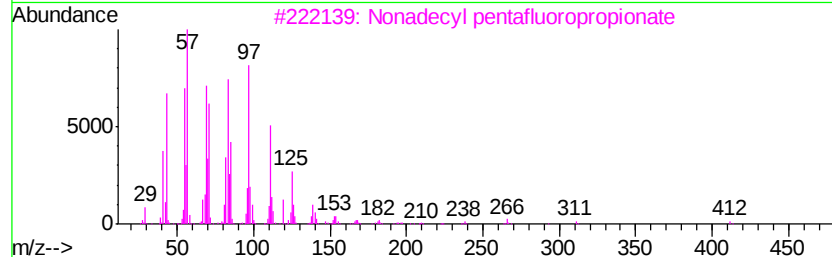
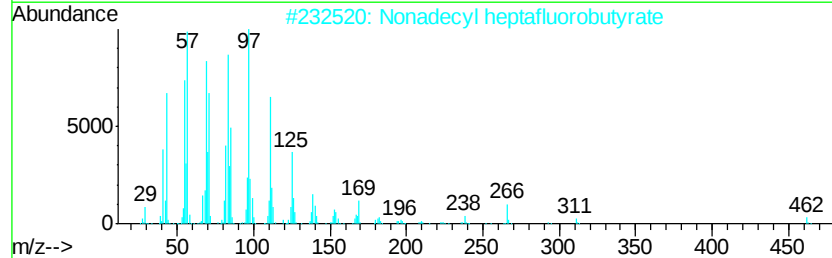
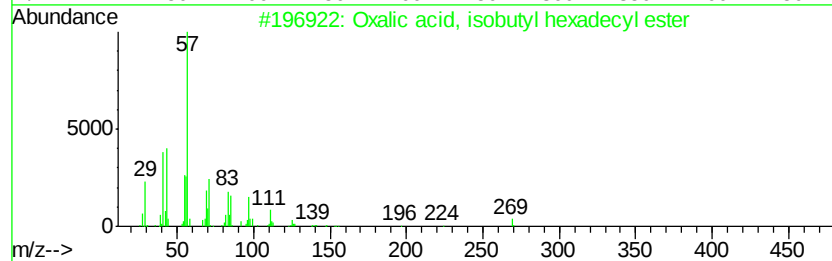
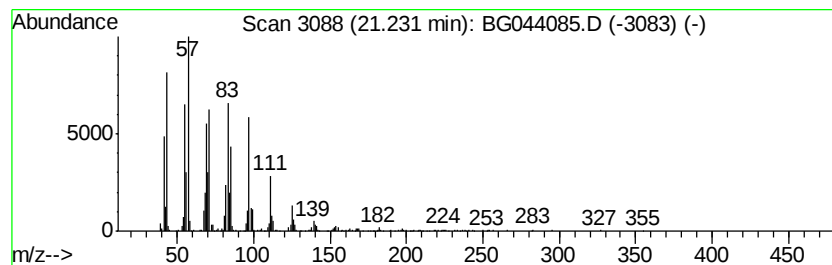
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Peak Number 6 Oxalic acid, isobutyl hexad... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.23	9.57 ng	620149	Chrysene-d12	21.56

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	91
2			Nonadecyl heptafluorobutyrate	480	C23H39F7O2	1000351-83-9	90
3			Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	87
4			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	87
5			Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	81



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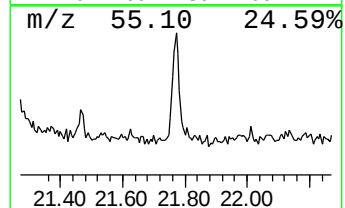
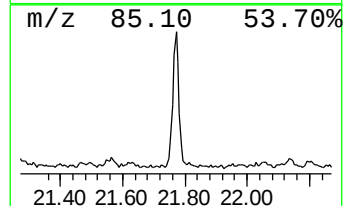
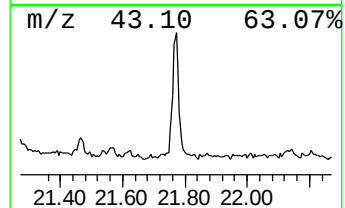
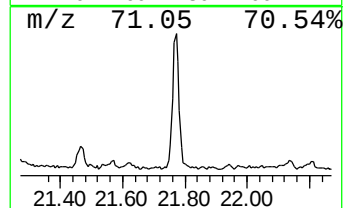
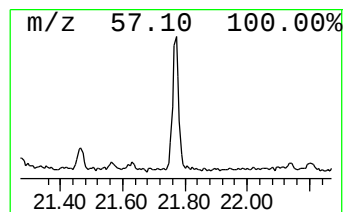
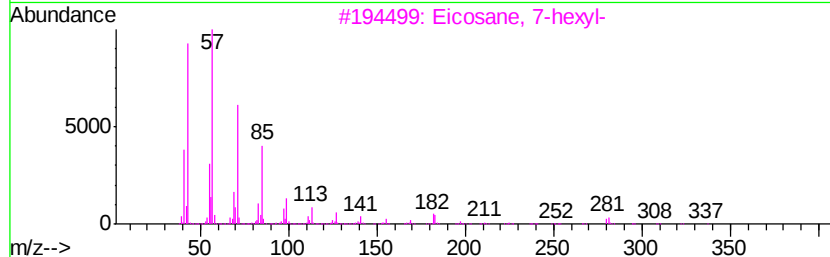
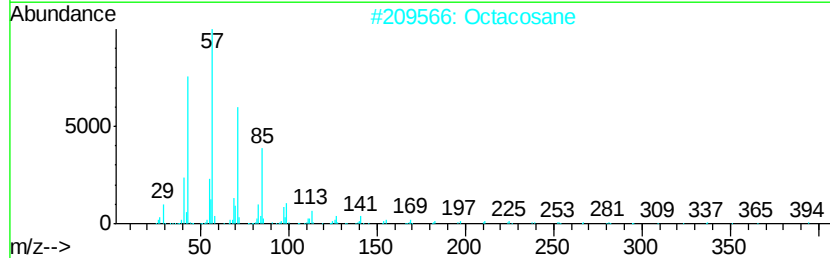
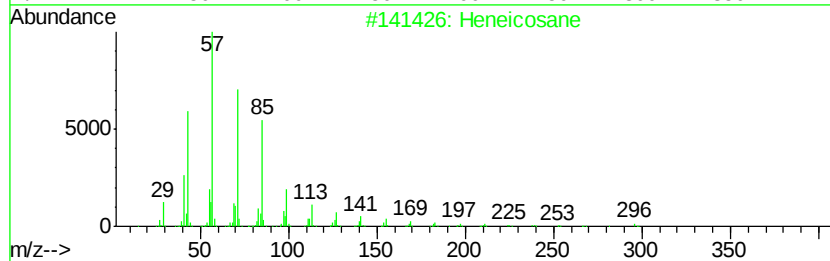
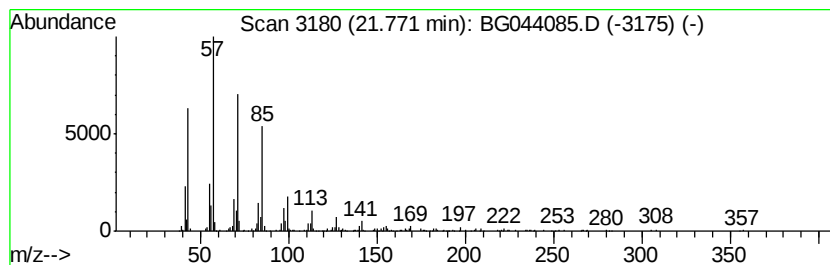
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heneicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.77	3.27 ng	211961	Chrysene-d12	21.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Octacosane		394	C28H58	000630-02-4	91
3	Eicosane, 7-hexyl-		366	C26H54	055333-99-8	91
4	Hexacosane		366	C26H54	000630-01-3	91
5	2-methyloctacosane		408	C29H60	1000376-72-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011720\
Data File : BG044085.D
Acq On : 17 Jan 2020 21:55
Operator : JU
Sample : L1162-09
Misc :
ALS Vial : 12 Sample Multiplier: 1

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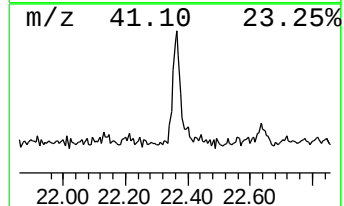
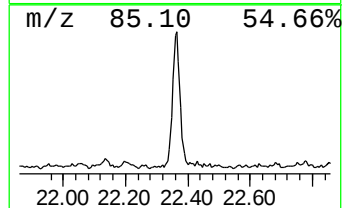
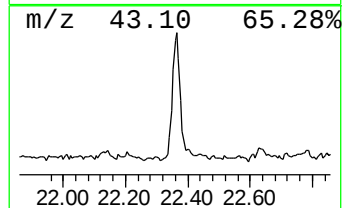
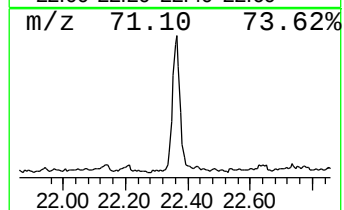
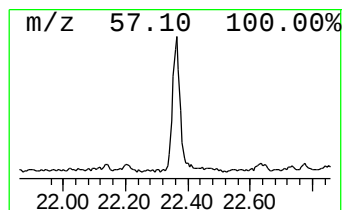
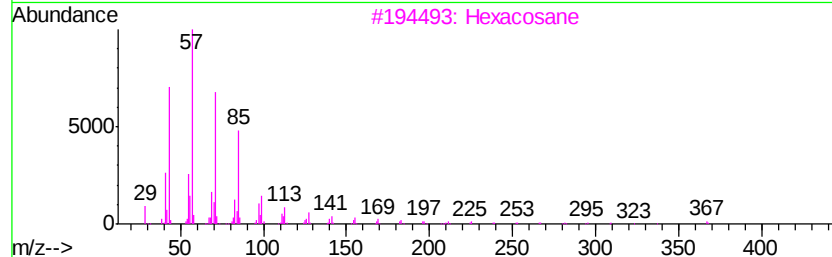
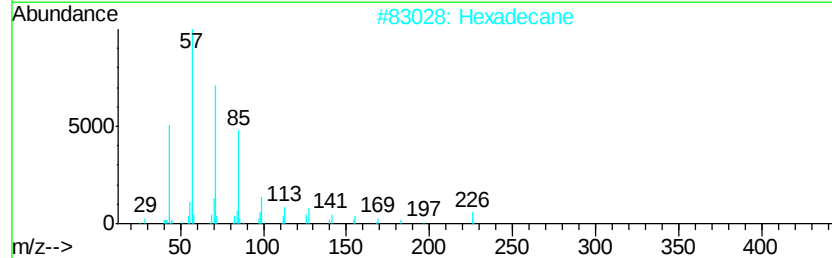
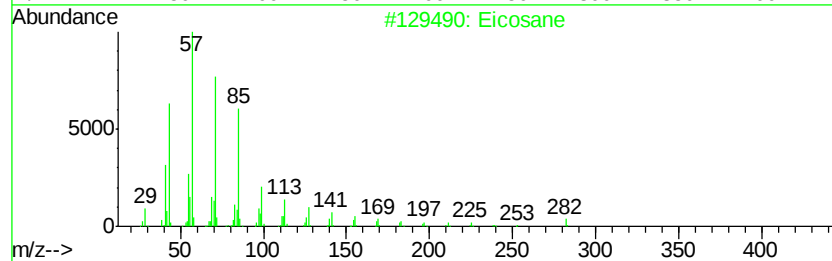
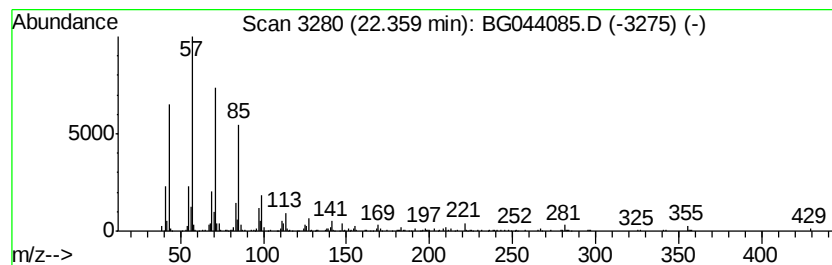
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Eicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.36	4.71 ng	304989	Chrysene-d12	21.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	96
2		Hexadecane	226	C16H34	000544-76-3	91
3		Hexacosane	366	C26H54	000630-01-3	90
4		Nonadecane	268	C19H40	000629-92-5	90
5		Octacosane	394	C28H58	000630-02-4	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011720\
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Misc :
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Instrument :
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ClientSampleId :
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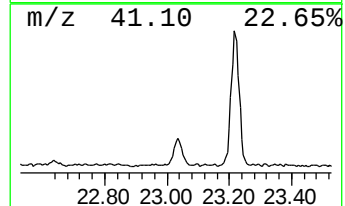
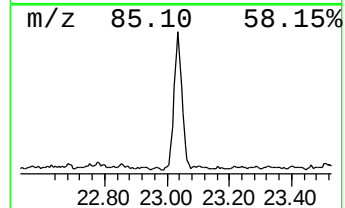
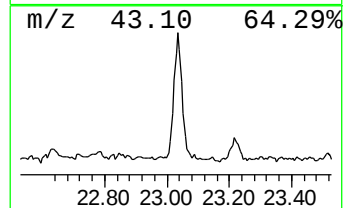
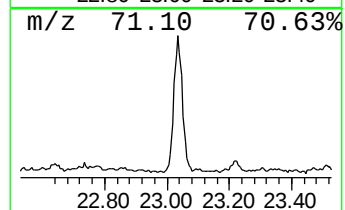
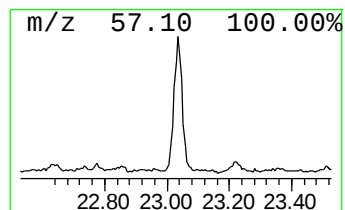
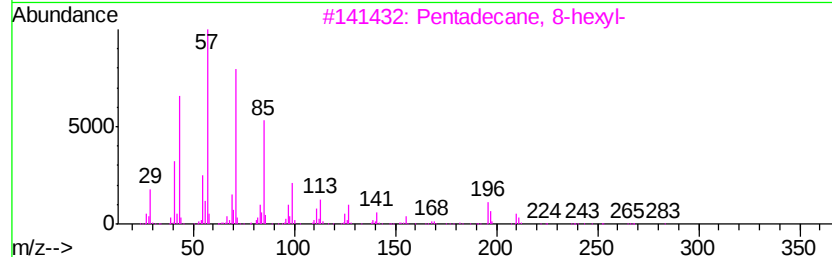
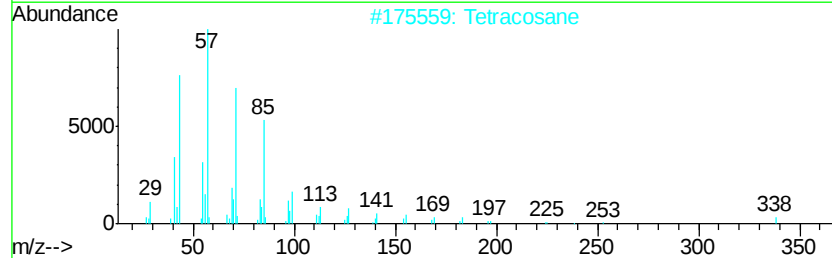
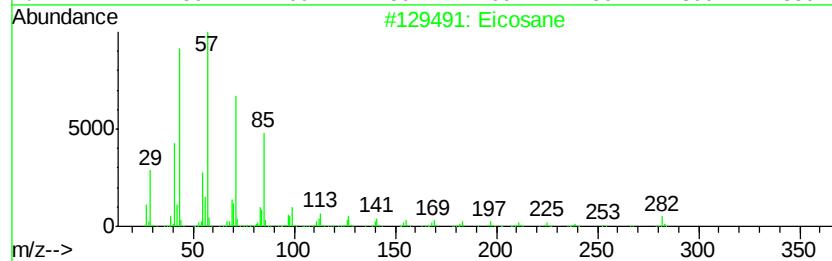
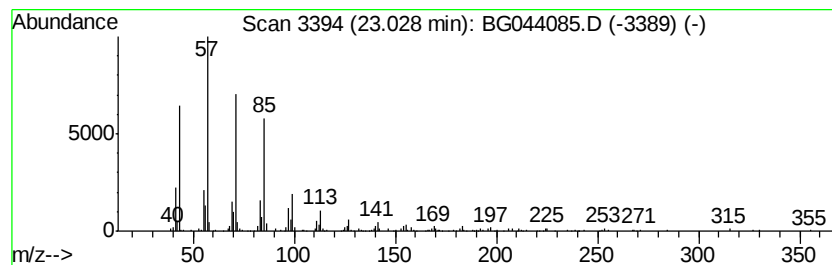
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tetracosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.03	5.06 ng	328209	Chrysene-d12	21.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	91
2	Tetracosane		338	C24H50	000646-31-1	91
3	Pentadecane, 8-hexyl-		296	C21H44	013475-75-7	91
4	Hexacosane		366	C26H54	000630-01-3	91
5	Eicosane, 7-hexyl-		366	C26H54	055333-99-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011720\
Data File : BG044085.D
Acq On : 17 Jan 2020 21:55
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Sample : L1162-09
Misc :
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Instrument :
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ClientSampleId :
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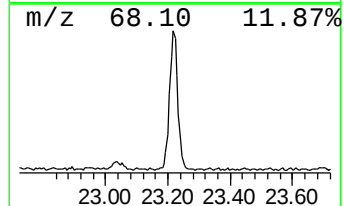
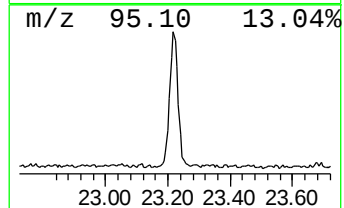
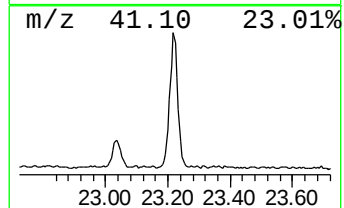
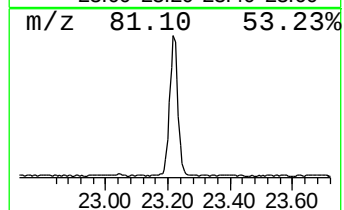
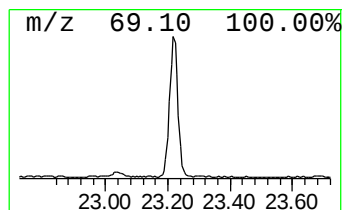
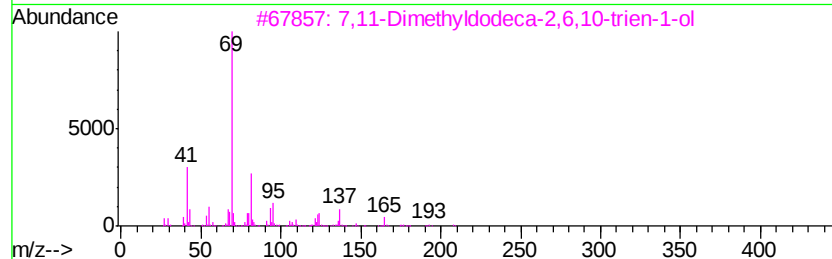
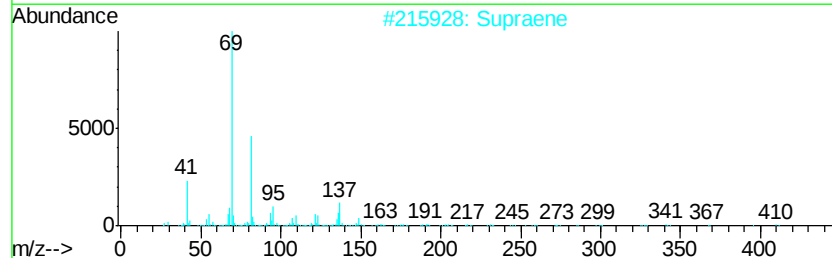
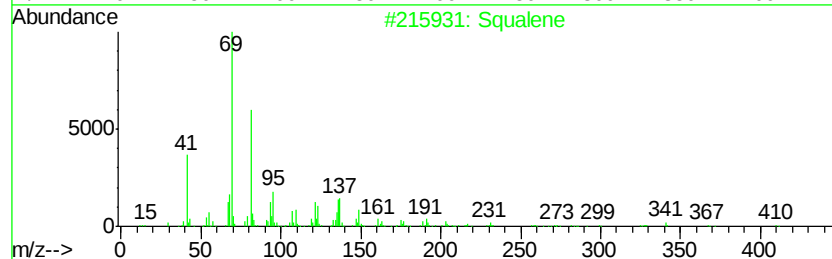
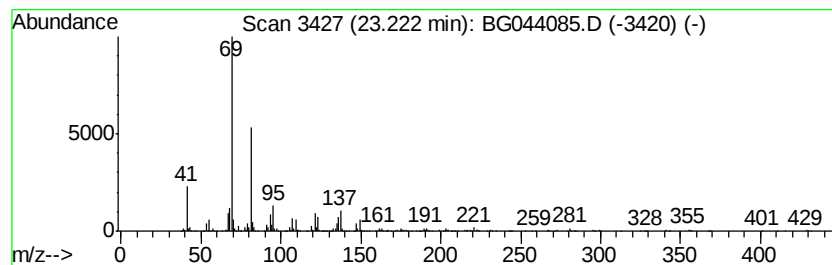
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Squalene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.22	15.40 ng	977334	Perylene-d12	24.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	000111-02-4	98
2		Supraene	410	C30H50	007683-64-9	90
3		7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	72
4		4,8,12-Tetradecatrienal, 5,9,13-...	248	C17H28O	066408-55-7	72
5		2,6,10,14-Hexadecatetraenoic aci...	332	C22H36O2	024035-35-6	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011720\
Data File : BG044085.D
Acq On : 17 Jan 2020 21:55
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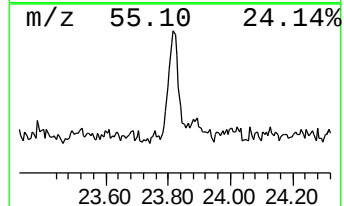
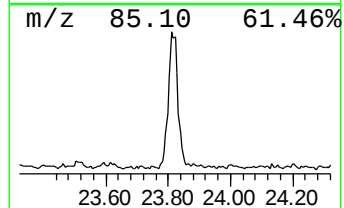
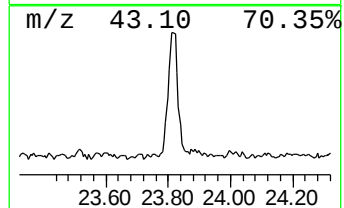
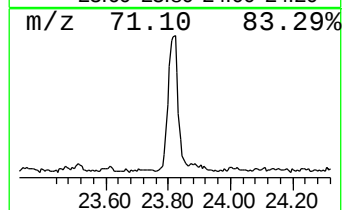
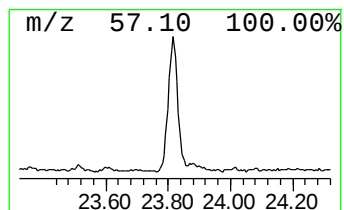
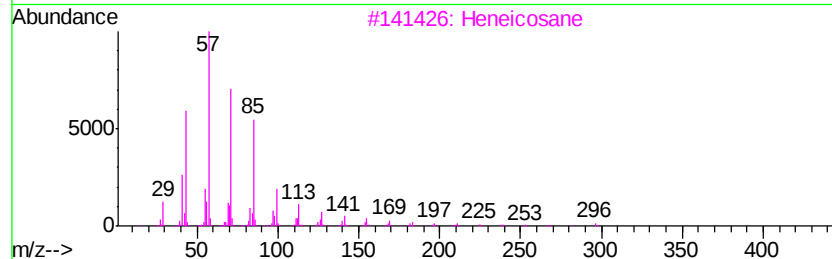
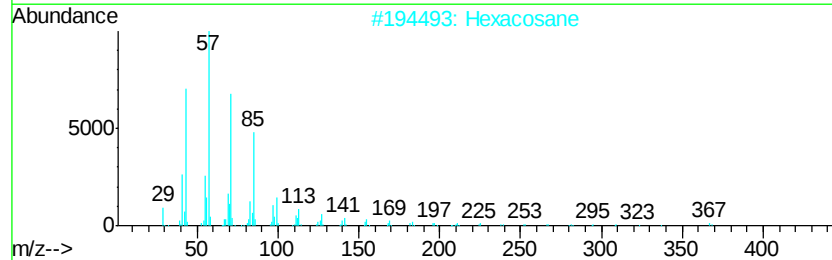
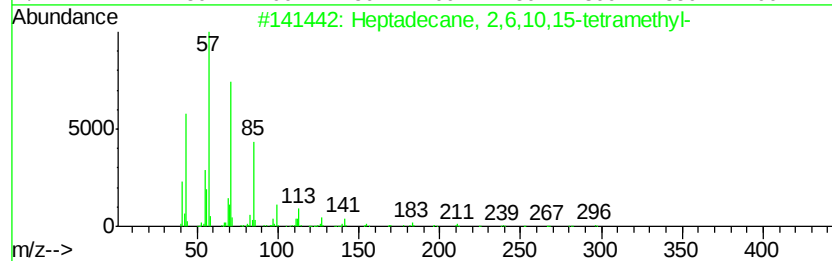
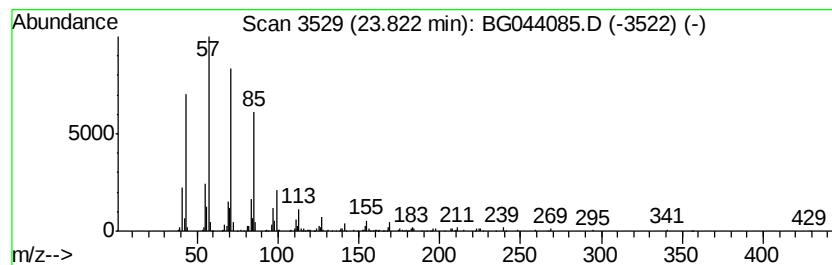
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Heptadecane, 2,6,10,15-tetr... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.82	4.27 ng	271221	Perylene-d12	24.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	93
2		Hexacosane	366	C26H54	000630-01-3	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Hexatriacontane	507	C36H74	000630-06-8	87
5		Eicosane	282	C20H42	000112-95-8	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011720\
 Data File : BG044085.D
 Acq On : 17 Jan 2020 21:55
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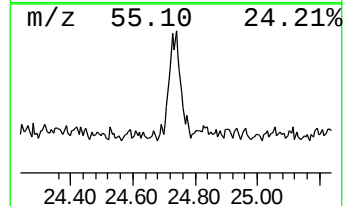
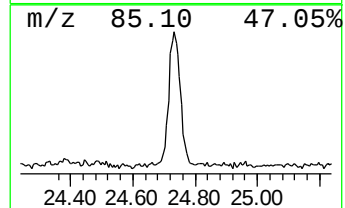
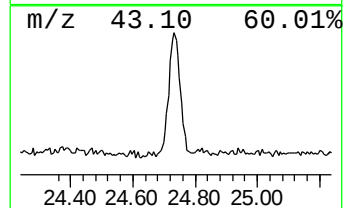
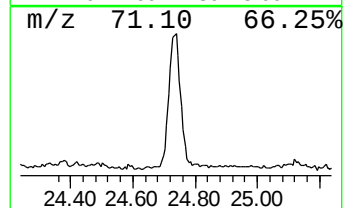
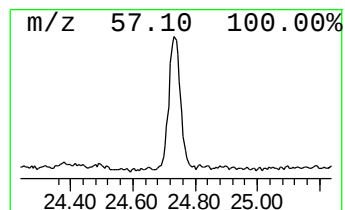
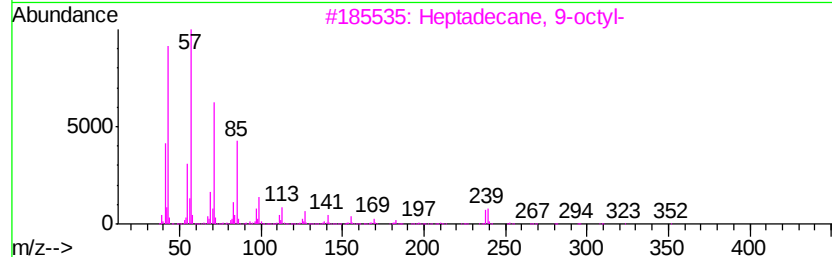
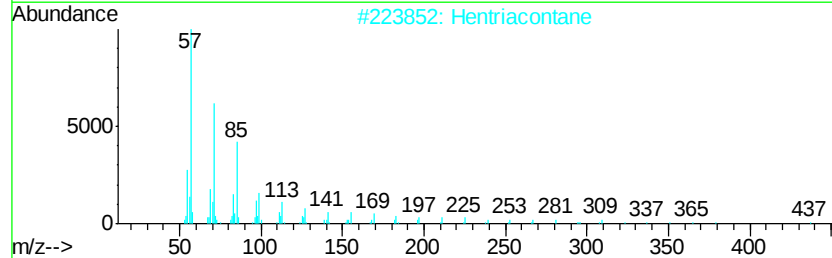
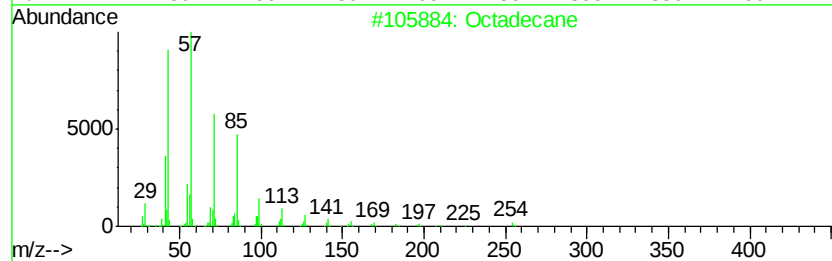
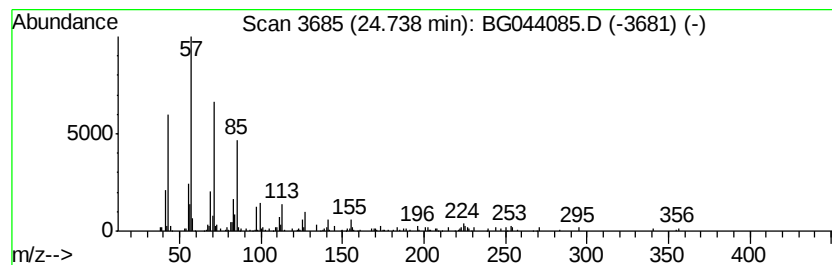
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 12 Octadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.74	3.66 ng	232581	Perylene-d12	24.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	93
2			Hentriacontane	437	C31H64	000630-04-6	87
3			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
4			Heneicosane, 11-decyl-	437	C31H64	055320-06-4	87
5			Tetracosane, 11-decyl-	479	C34H70	055429-84-0	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011720\
Data File : BG044085.D
Acq On : 17 Jan 2020 21:55
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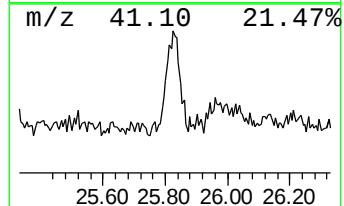
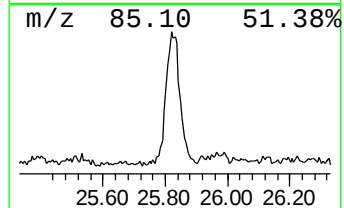
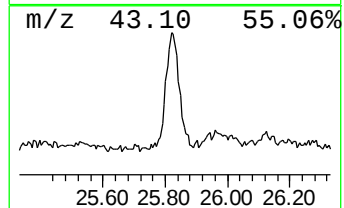
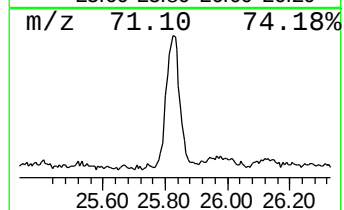
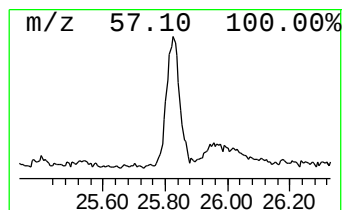
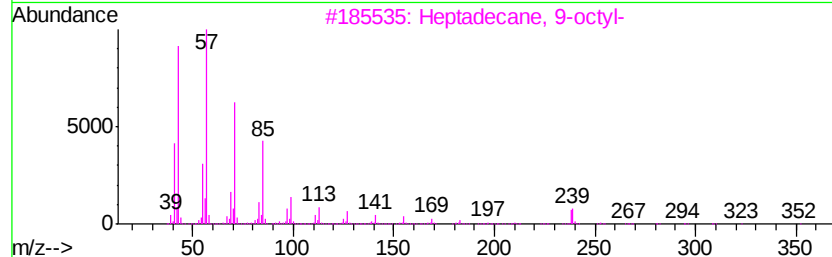
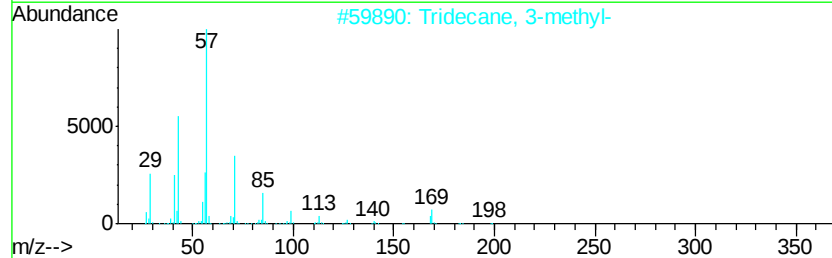
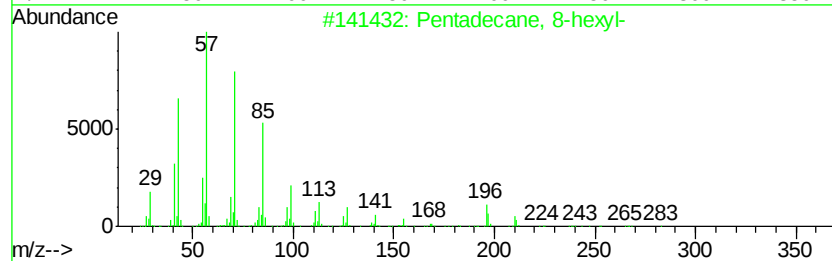
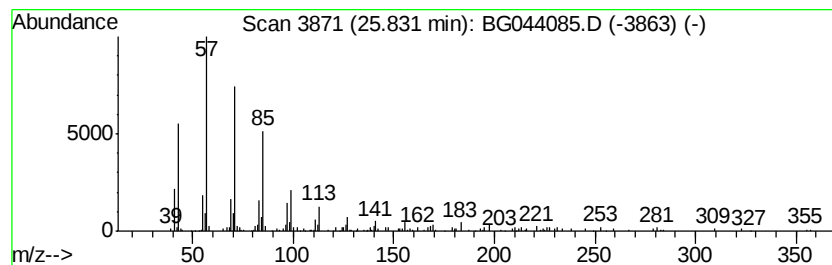
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Pentadecane, 8-hexyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.83	3.25 ng	206124	Perylene-d12	24.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	87
2		Tridecane, 3-methyl-	198	C14H30	006418-41-3	87
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
4		Dodecane, 2-methyl-	184	C13H28	001560-97-0	87
5		6,6-Diethylhoctadecane	310	C22H46	1000360-41-8	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG011720\
 Data File : BG044085.D
 Acq On : 17 Jan 2020 21:55
 Operator : JU
 Sample : L1162-09
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MH-719

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
2-Pentanone, 4-hy...	5.04	7.3	ng	178818	1	7.93	487007	20.0
unknown7.46	7.46	101.4	ng	2468090	1	7.93	487007	20.0
.alpha.-Ethyl-.al...	9.03	2.4	ng	59525	1	7.93	487007	20.0
2-Naphthalenamine...	20.23	2.4	ng	153204	5	21.56	1296240	20.0
Oxalic acid, isob...	21.23	9.6	ng	620149	5	21.56	1296240	20.0
Heneicosane	21.77	3.3	ng	211961	5	21.56	1296240	20.0
Eicosane	22.36	4.7	ng	304989	5	21.56	1296240	20.0
Tetracosane	23.03	5.1	ng	328209	5	21.56	1296240	20.0
Squalene	23.22	15.4	ng	977334	6	24.67	1269230	20.0
Heptadecane, 2,6,...	23.82	4.3	ng	271221	6	24.67	1269230	20.0
Octadecane	24.74	3.7	ng	232581	6	24.67	1269230	20.0
Pentadecane, 8-he...	25.83	3.3	ng	206124	6	24.67	1269230	20.0