

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011420\
 Data File : BG044036.D
 Acq On : 14 Jan 2020 14:49
 Operator : JU
 Sample : L1108-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 INWOOD-BH-01-A

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	3.135	3	8	30	rVB	516652	1238763	27.54%	2.806%
2	5.057	328	335	345	rVB	330183	538343	11.97%	1.220%
3	5.527	407	415	428	rVB	1446799	2353573	52.32%	5.332%
4	7.113	677	685	697	rBV	1496060	2671105	59.37%	6.051%
5	7.483	740	748	757	rBV	1515623	2629494	58.45%	5.957%
6	7.947	820	827	835	rBV	413226	710196	15.79%	1.609%
7	8.271	871	882	890	rBV	1150999	2022797	44.96%	4.583%
8	9.099	1014	1023	1033	rBV	882385	1596912	35.50%	3.618%
9	10.744	1294	1303	1315	rBV	547771	1003432	22.30%	2.273%
10	13.200	1714	1721	1729	rBV	2109418	3353113	74.53%	7.597%
11	14.029	1855	1862	1868	rBV	138417	213303	4.74%	0.483%
12	14.575	1948	1955	1962	rBV	860596	1356122	30.14%	3.072%
13	16.061	2201	2208	2221	rBV	1831522	2811920	62.50%	6.370%
14	17.319	2415	2422	2426	rBV	1028044	1578921	35.10%	3.577%
15	17.360	2426	2429	2436	rVV	194909	289109	6.43%	0.655%
16	17.448	2440	2444	2449	rVB	36668	58960	1.31%	0.134%
17	18.224	2568	2576	2586	rBV2	120355	283202	6.30%	0.642%
18	18.394	2600	2605	2608	rBV3	30607	52669	1.17%	0.119%
19	19.369	2766	2771	2777	rBV	291750	421718	9.37%	0.955%
20	19.728	2824	2832	2839	rBV	234400	400072	8.89%	0.906%
21	19.933	2862	2867	2878	rVB	3214184	4498811	100.00%	10.192%
22	20.656	2971	2990	3004	rBV	858131	4219743	93.80%	9.560%
23	21.238	3084	3089	3096	rBV2	507548	794117	17.65%	1.799%
24	21.302	3096	3100	3108	rVB2	133290	211452	4.70%	0.479%
25	21.473	3125	3129	3135	rBV	156515	224214	4.98%	0.508%
26	21.573	3138	3146	3151	rBV	985704	1831834	40.72%	4.150%
27	21.784	3179	3182	3208	rVB2	352621	1289290	28.66%	2.921%
28	22.384	3278	3284	3293	rBV2	262303	494810	11.00%	1.121%
29	22.536	3297	3310	3329	rVB4	231778	1391913	30.94%	3.153%
30	23.059	3394	3399	3412	rVB	247311	477957	10.62%	1.083%
31	23.241	3425	3430	3444	rVB	255662	626976	13.94%	1.420%
32	23.852	3528	3534	3539	rBV	197767	438045	9.74%	0.992%
33	24.698	3668	3678	3686	rBV2	557429	1772709	39.40%	4.016%
34	25.897	3876	3882	3891	rVB2	107483	284402	6.32%	0.644%

LSC Area Percent Report

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Sample : L1108-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
INWOOD-BH-01-A

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

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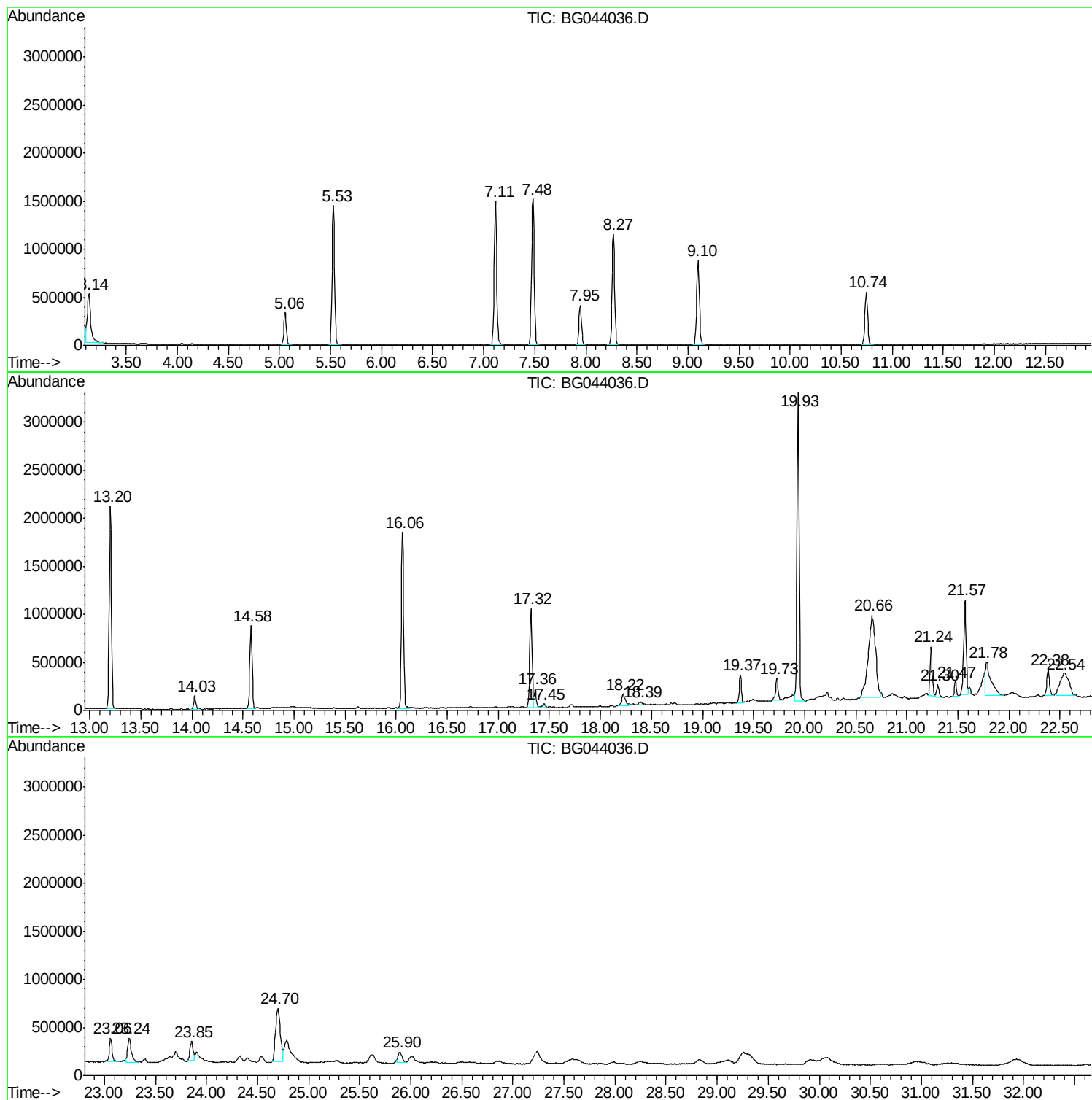
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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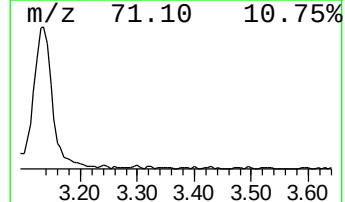
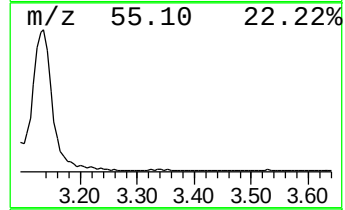
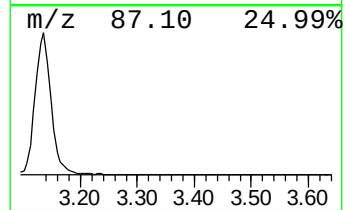
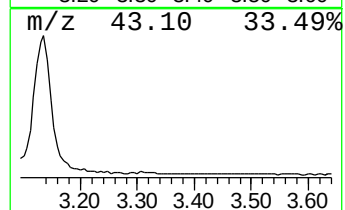
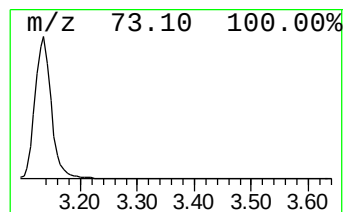
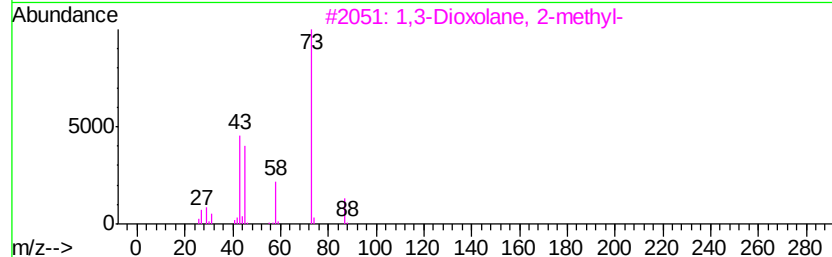
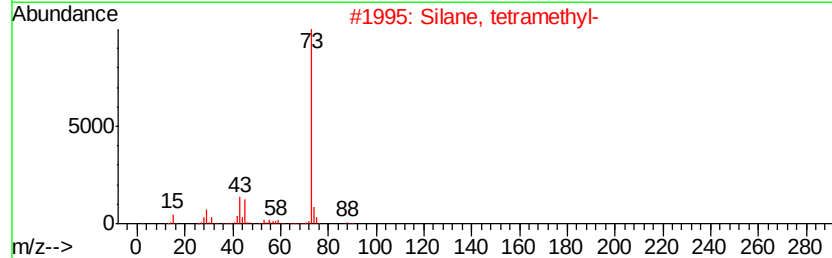
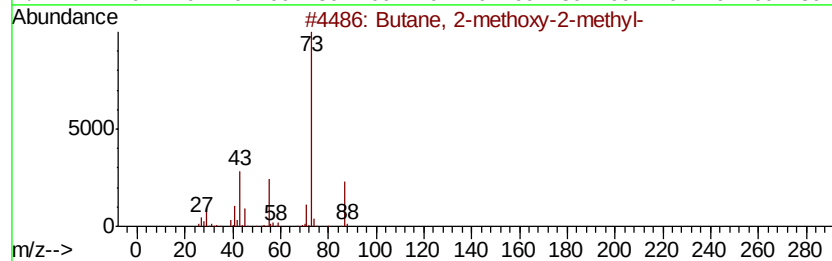
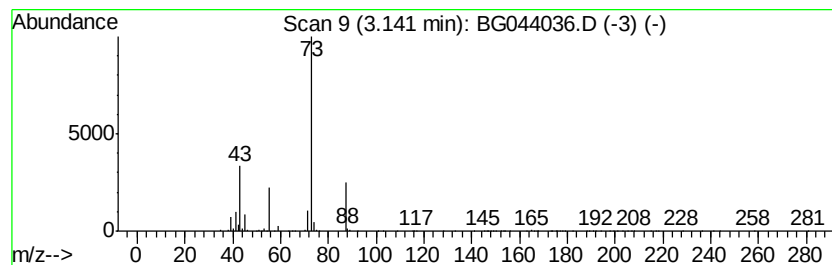
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.14	34.89 ng	1238760	1,4-Dichlorobenzene-d4	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
2		Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
4		N-Ethylformamide	73	C3H7NO	000627-45-2	9
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9



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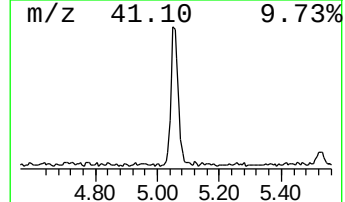
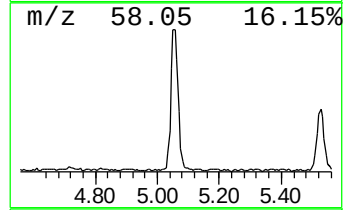
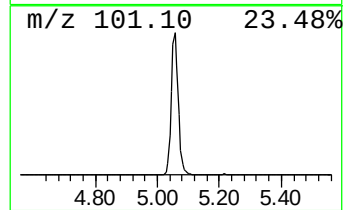
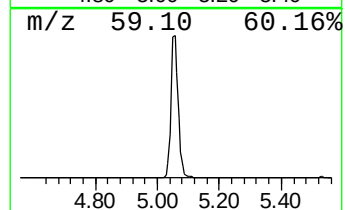
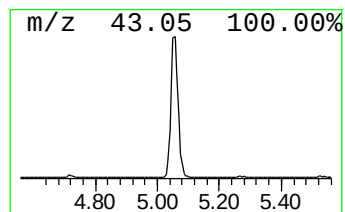
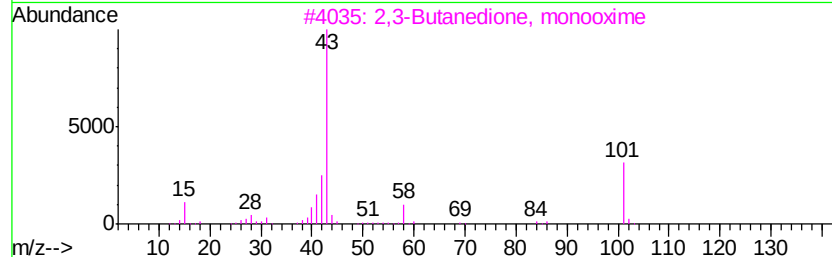
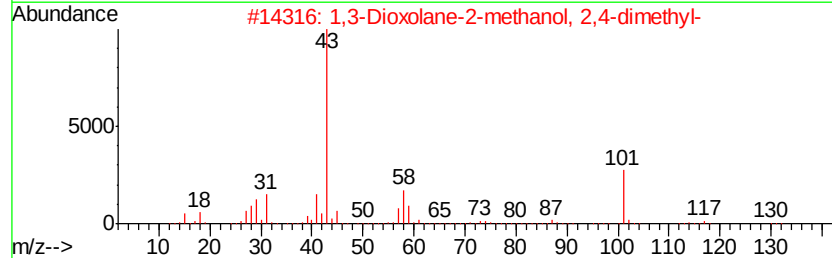
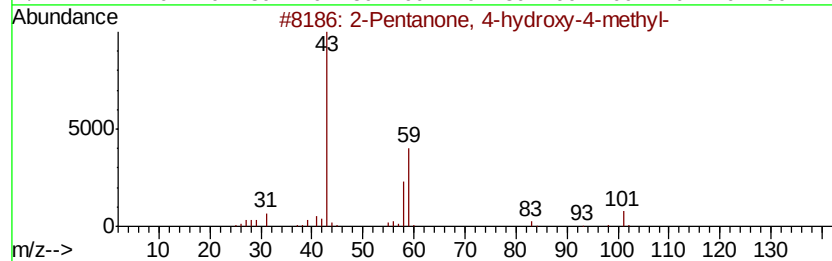
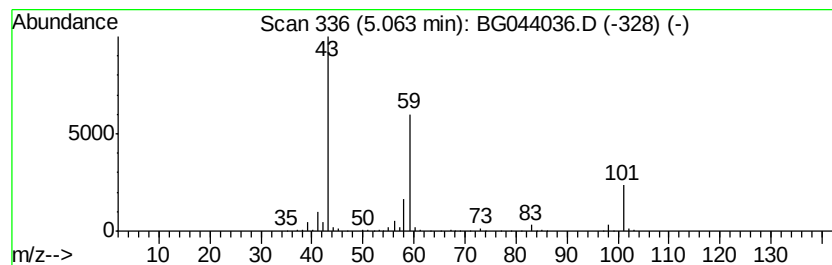
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.06	15.16 ng	538343	1,4-Dichlorobenzene-d4	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		1,3-Dioxolane-2-methanol, 2,4-di...	132	C6H12O3	053951-43-2	28
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4		4-Hydroxy-3-hexanone	116	C6H12O2	004984-85-4	23
5		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17



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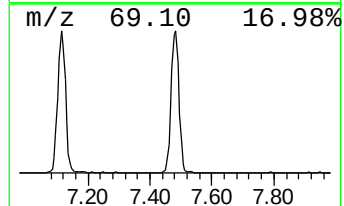
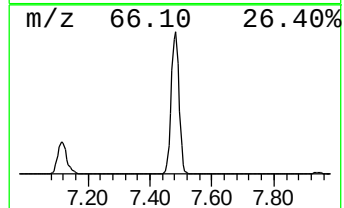
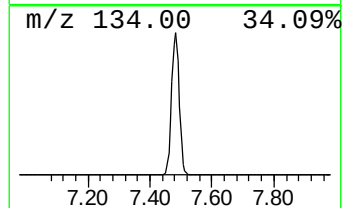
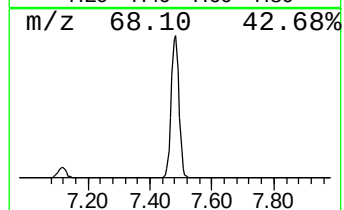
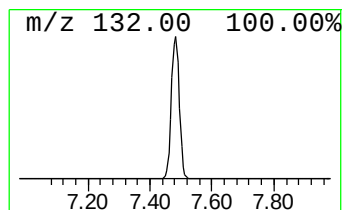
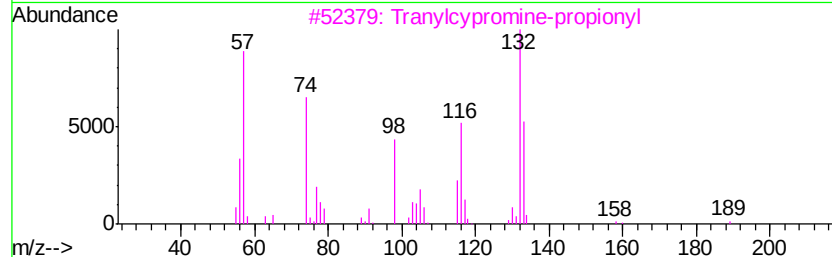
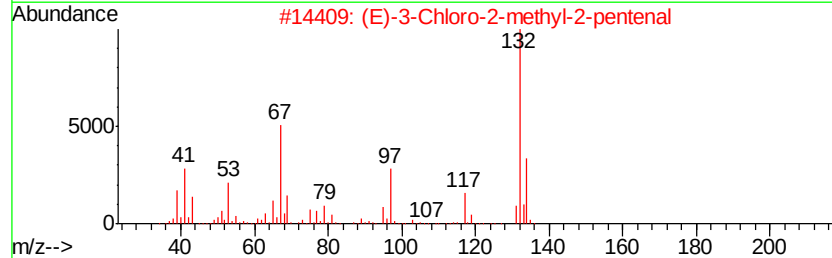
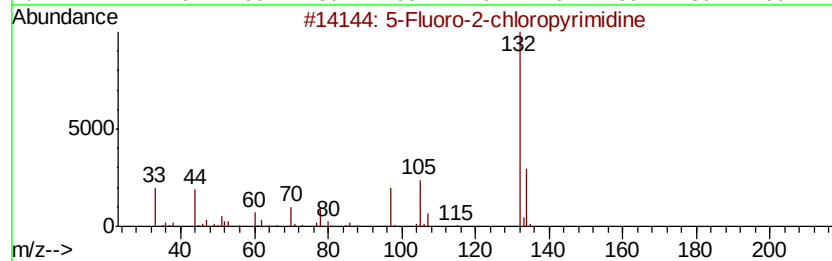
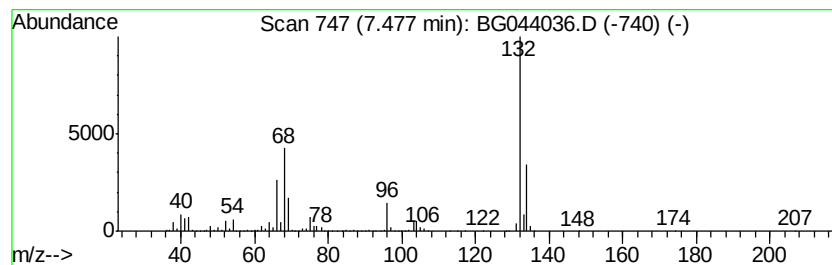
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown7.48 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.48	74.05 ng	2629490	1,4-Dichlorobenzene-d4	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		Tranvlcypramine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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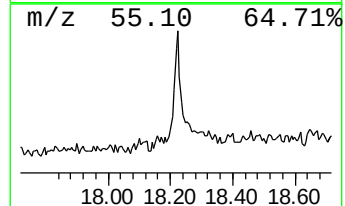
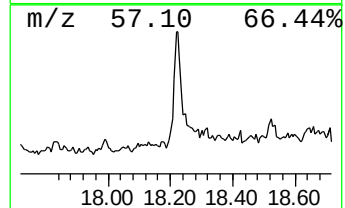
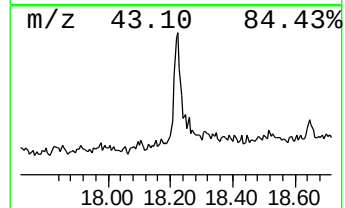
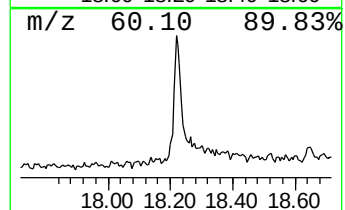
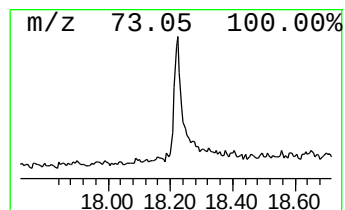
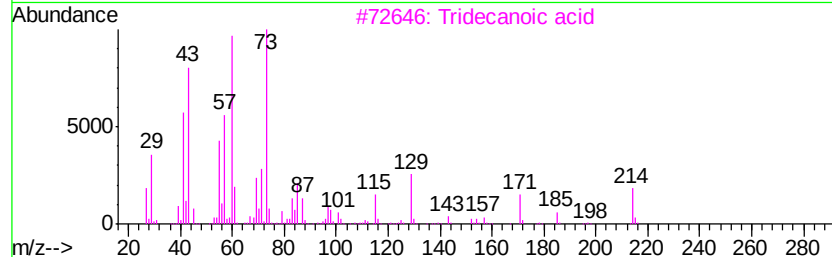
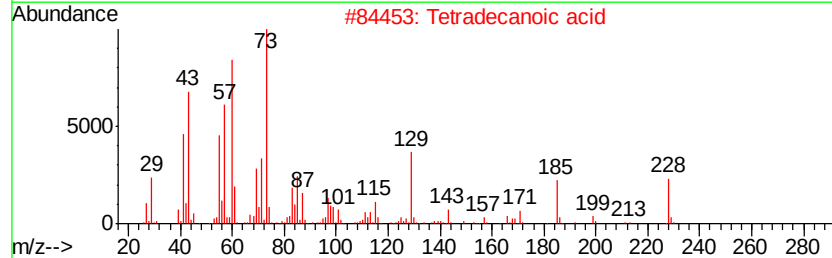
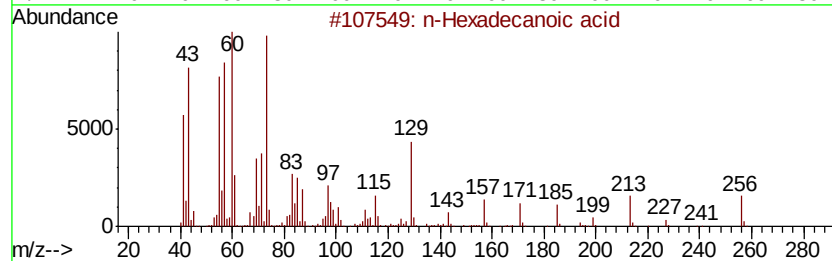
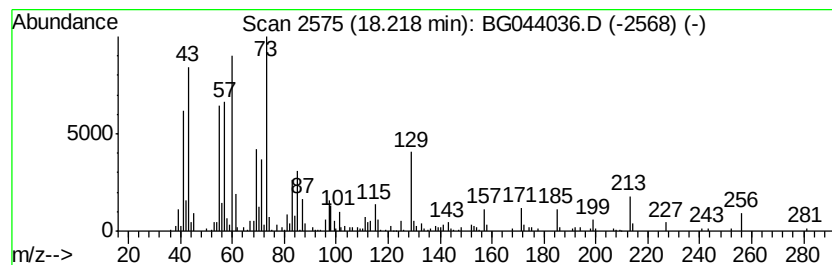
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 n-Hexadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.22	3.59 ng	283202	Phenanthrene-d10	17.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
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	64



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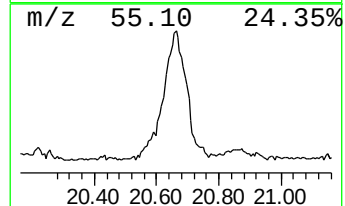
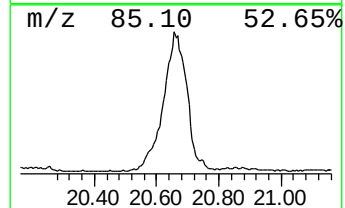
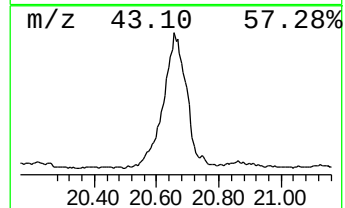
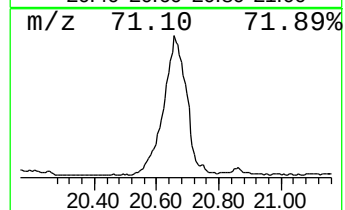
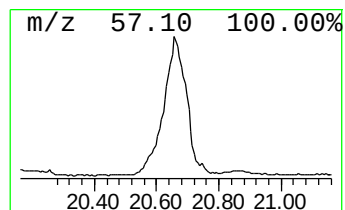
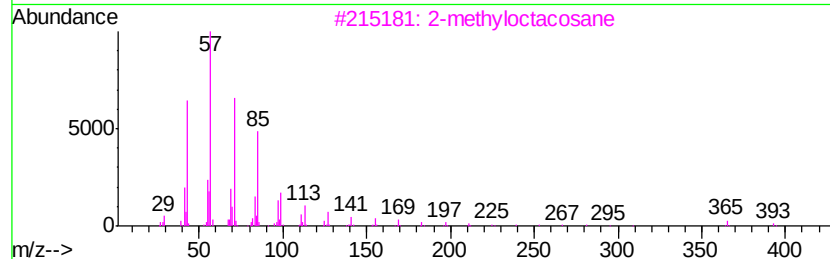
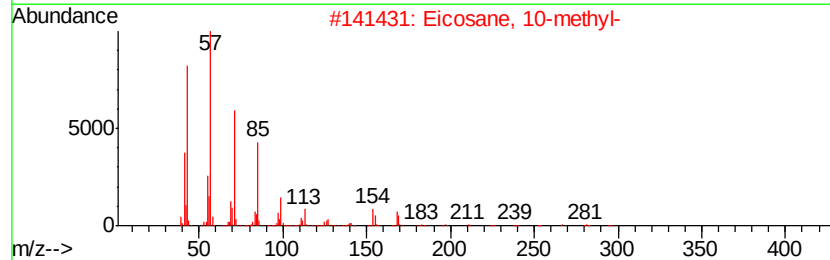
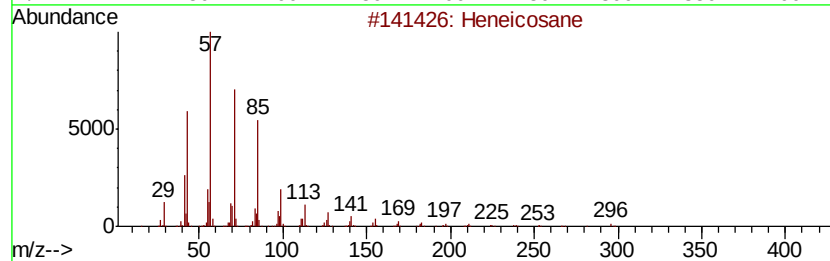
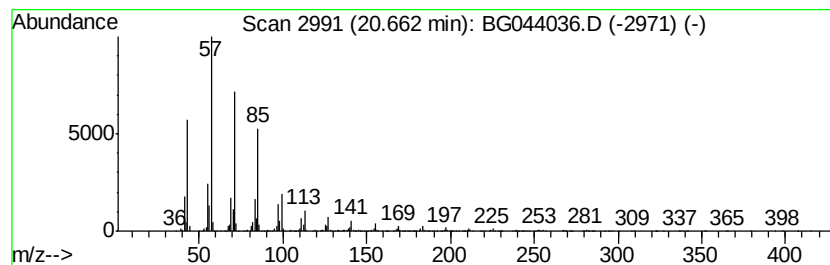
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Peak Number 6 Heneicosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.66	46.07 ng	4219740	Chrysene-d12	21.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	94
2		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Octacosane	394	C28H58	000630-02-4	91
5		Hexacosane	366	C26H54	000630-01-3	91



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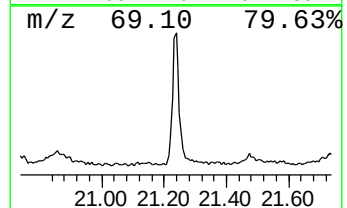
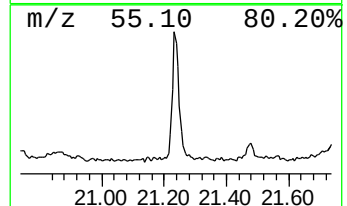
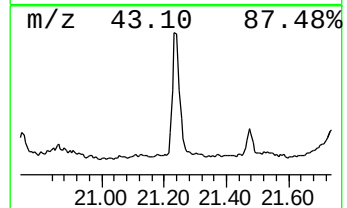
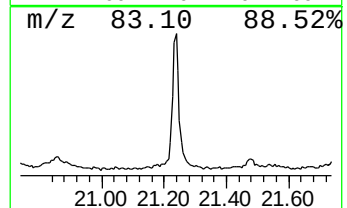
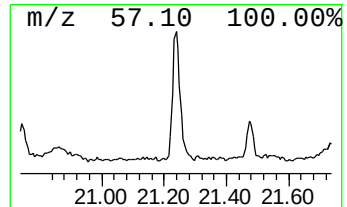
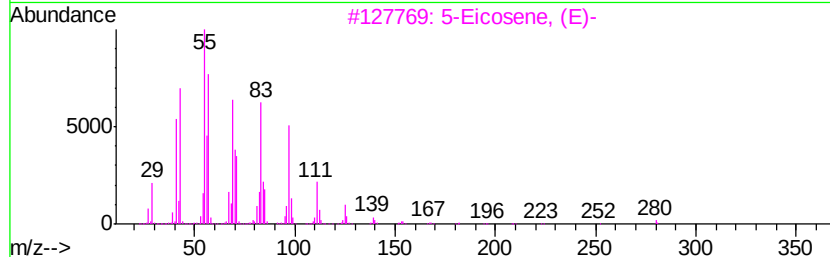
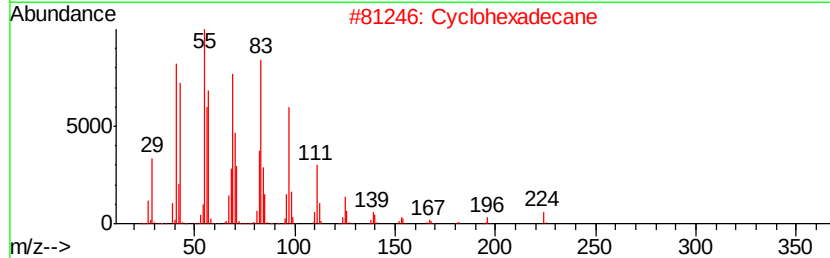
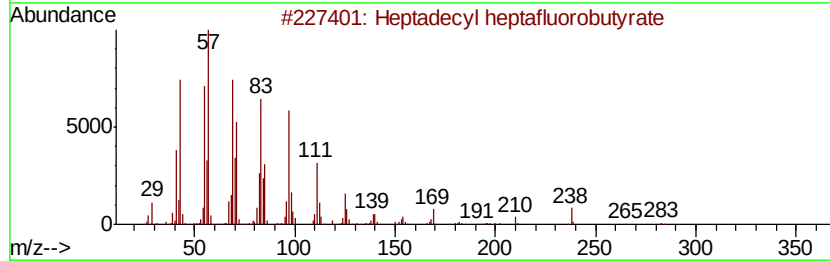
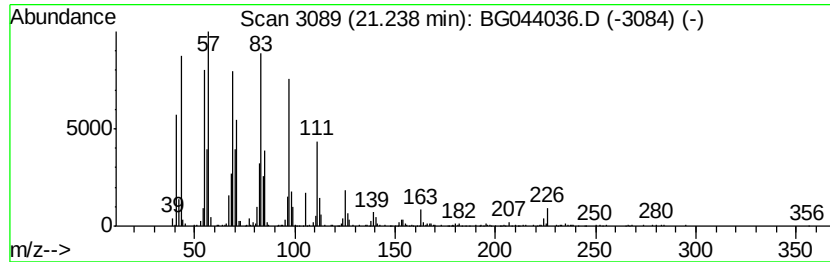
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Heptadecyl heptafluorobutyrate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.24	8.67 ng	794117	Chrysene-d12	21.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
2		Cyclohexadecane	224	C16H32	000295-65-8	93
3		5-Eicosene, (E)-	280	C20H40	074685-30-6	93
4		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	93
5		1-Heneicosanol	312	C21H44O	015594-90-8	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011420\
Data File : BG044036.D
Acq On : 14 Jan 2020 14:49
Operator : JU
Sample : L1108-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

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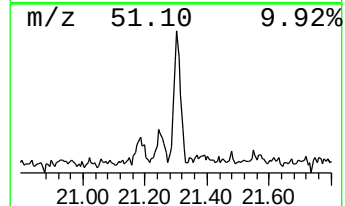
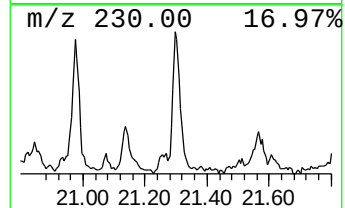
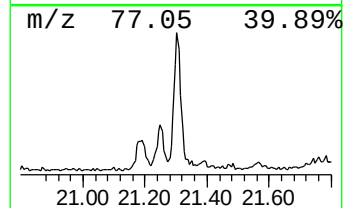
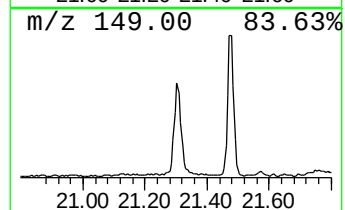
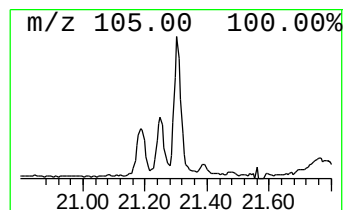
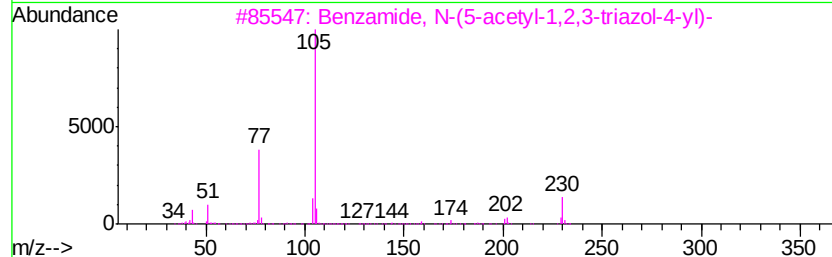
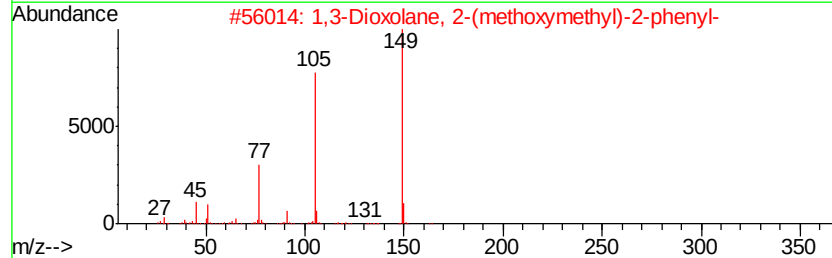
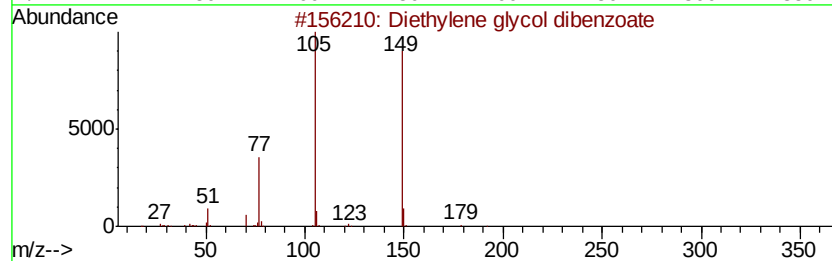
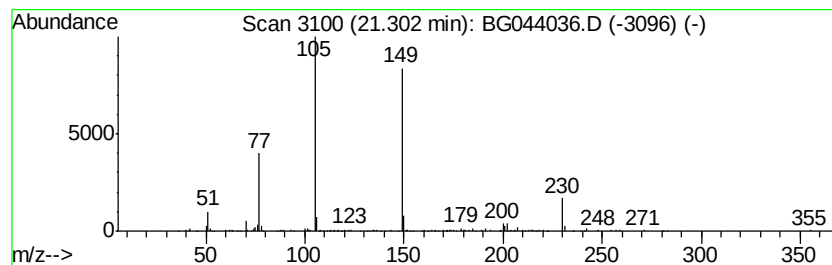
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Diethylene glycol dibenzoate Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.30	2.31 ng	211452	Chrysene-d12	21.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	72
2		1,3-Dioxolane, 2-(methoxymethyl)...	194	C11H14O3	1000156-71-2	59
3		Benzamide, N-(5-acetyl-1,2,3-tri...	230	C11H10N4O2	129959-33-7	50
4		2-(2-Carboxyvinyl)pyridine, trans	149	C8H7NO2	007340-22-9	49
5		2,2'-(Ethane-1,2-diylbis(oxy))bi...	358	C20H22O6	1000366-99-4	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011420\
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Misc :
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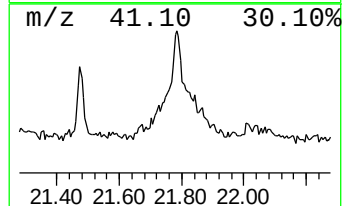
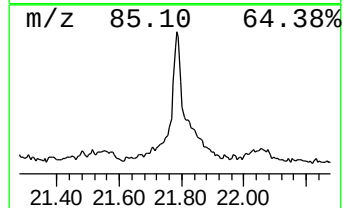
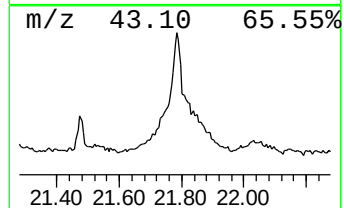
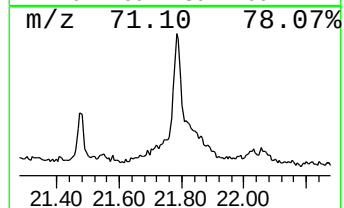
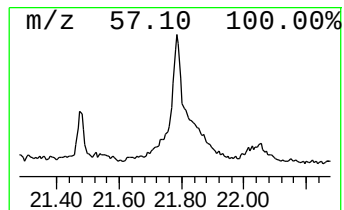
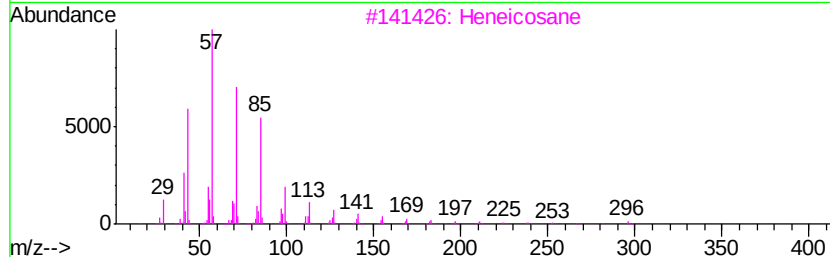
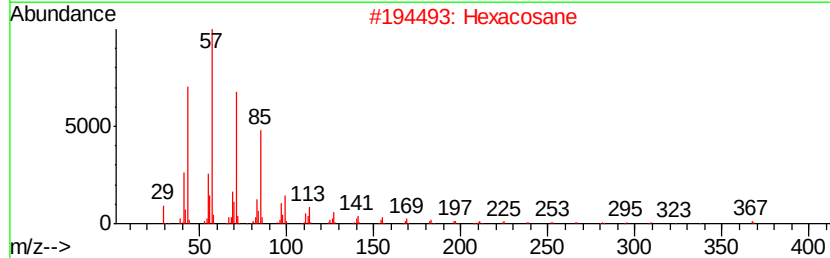
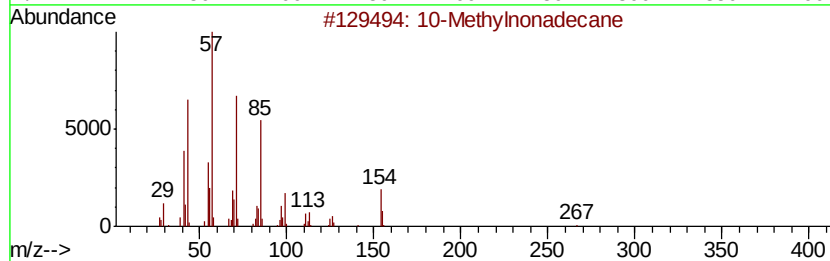
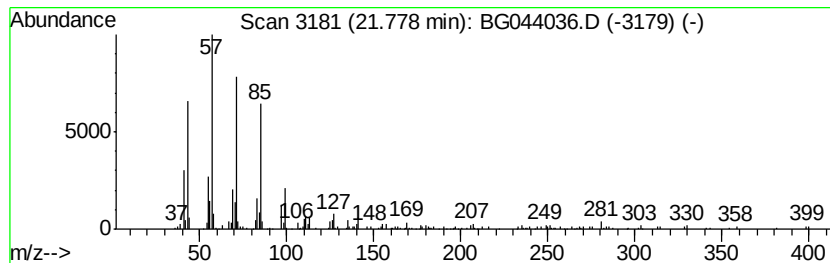
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 10-Methylnonadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.78	14.08 ng	1289290	Chrysene-d12	21.57

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	10-Methyl	nonadecane	282	C20H42	056862-62-5	87	
2	Hexacosane	366	C26H54	000630-01-3	87		
3	Heneicosane	296	C21H44	000629-94-7	86		
4	Octacosane	394	C28H58	000630-02-4	80		
5	Octadecane	254	C18H38	000593-45-3	72		



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011420\
Data File : BG044036.D
Acq On : 14 Jan 2020 14:49
Operator : JU
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ALS Vial : 5 Sample Multiplier: 1

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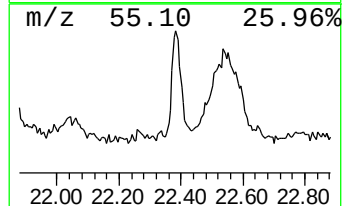
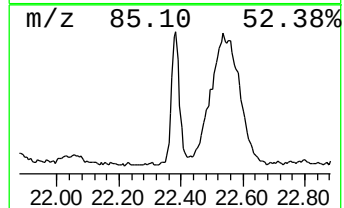
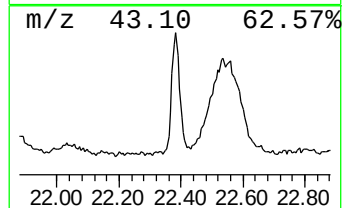
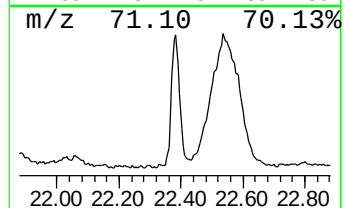
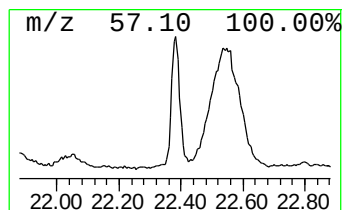
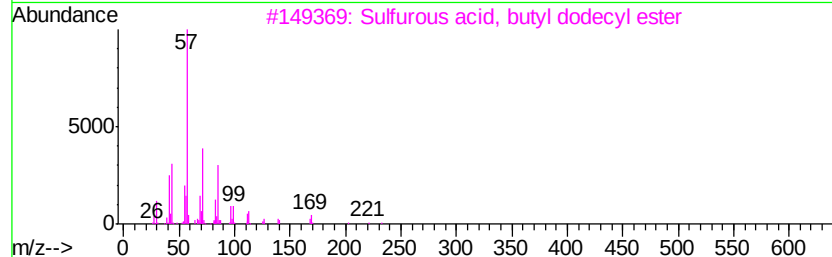
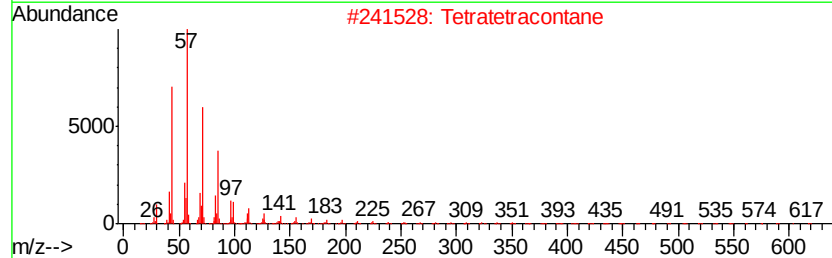
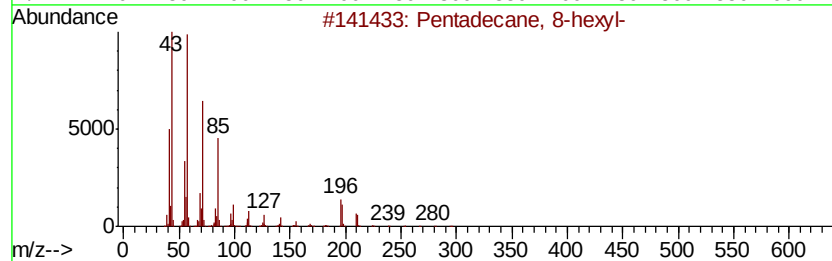
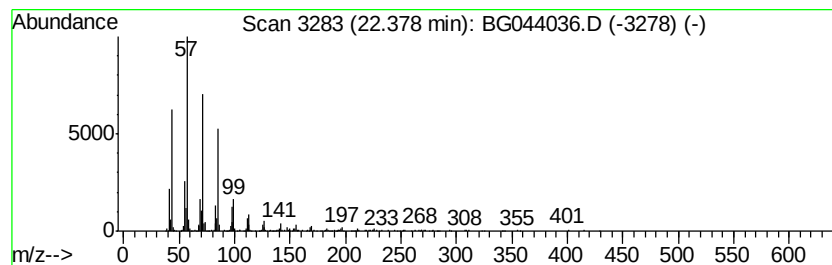
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Pentadecane, 8-hexyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.38	5.40 ng	494810	Chrysene-d12	21.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	90
2		Tetratetracontane	619	C44H90	007098-22-8	90
3		Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	87
4		Octacosane	394	C28H58	000630-02-4	87
5		10-Methylnonadecane	282	C20H42	056862-62-5	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011420\
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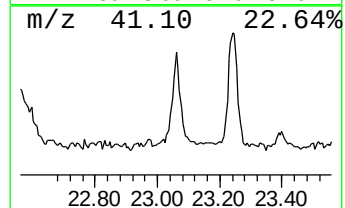
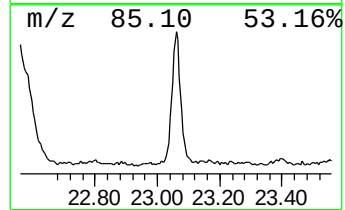
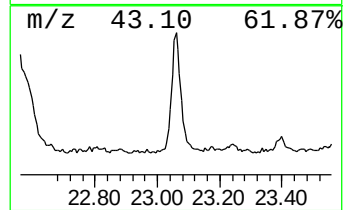
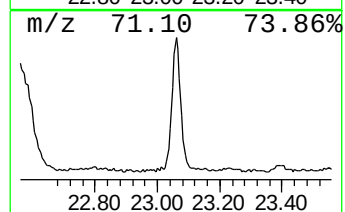
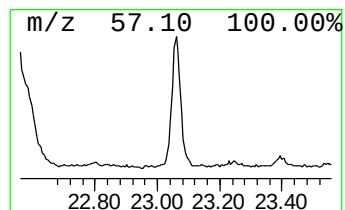
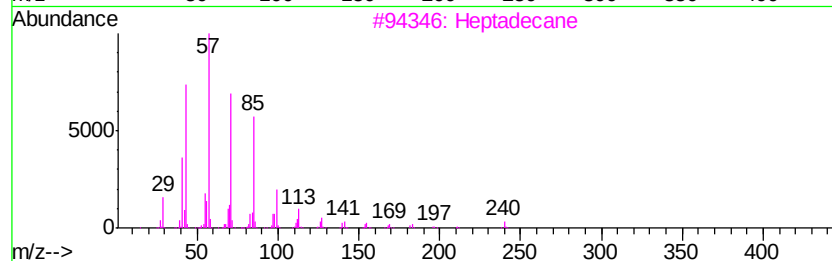
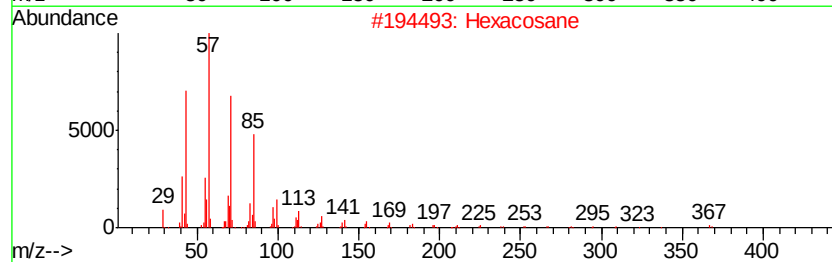
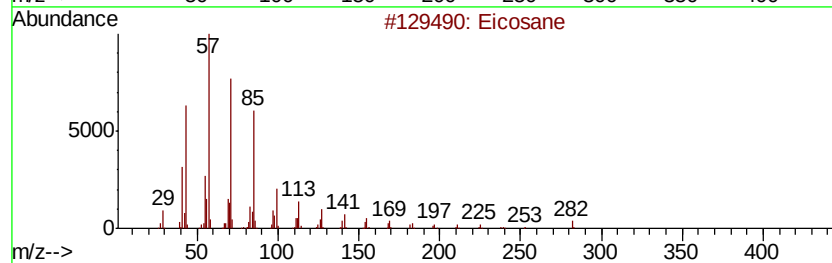
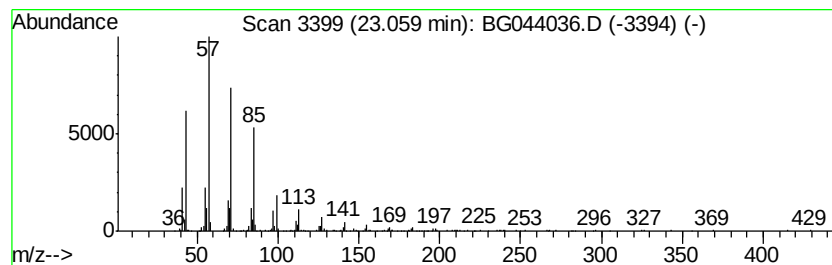
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TIC Library : C:\Database\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 12 Eicosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.06	5.22 ng	477957	Chrysene-d12	21.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	94
2		Hexacosane	366	C26H54	000630-01-3	91
3		Heptadecane	240	C17H36	000629-78-7	91
4		Nonadecane	268	C19H40	000629-92-5	90
5		Tetracosane	338	C24H50	000646-31-1	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011420\
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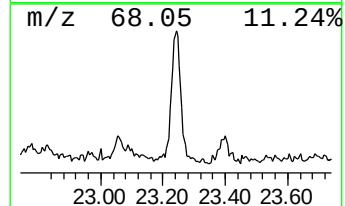
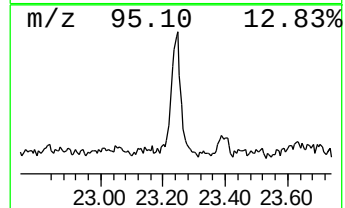
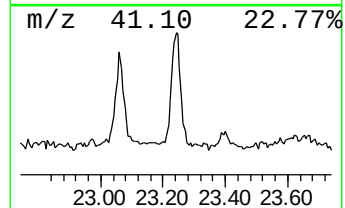
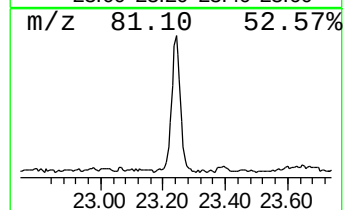
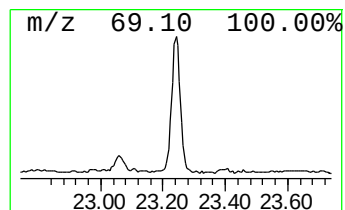
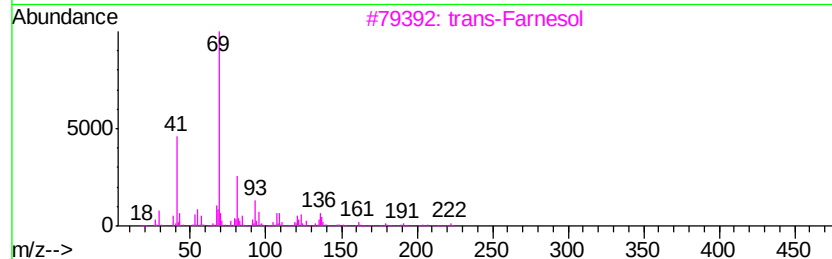
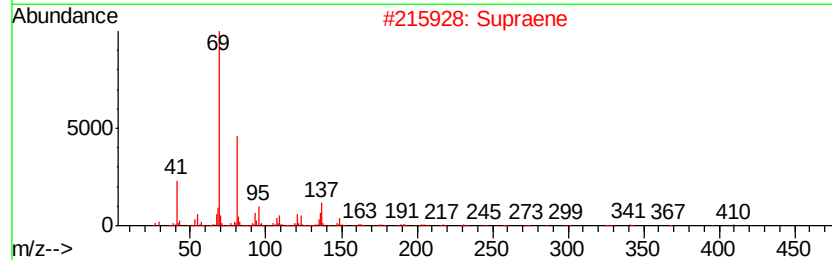
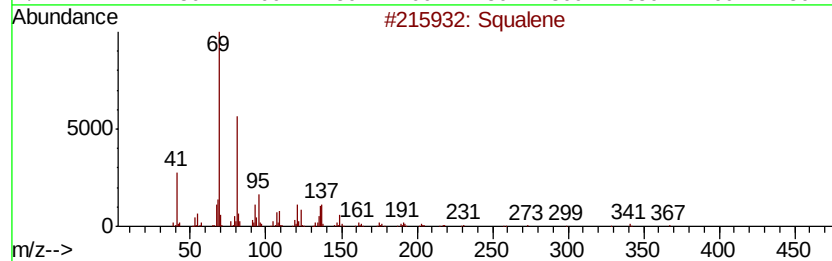
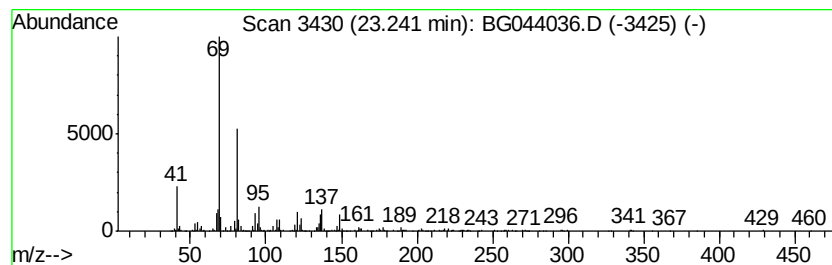
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Squalene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.24	7.07 ng	626976	Perylene-d12	24.70
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Squalene	410 C30H50	000111-02-4	91
2	Supraene	410 C30H50	007683-64-9	91
3	trans-Farnesol	222 C15H26O	000106-28-5	64
4	Methanone, dicyclopropyl-	110 C7H10O	001121-37-5	47
5	2,6-Octadiene, 4,5-dimethyl-	138 C10H18	018476-57-8	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011420\
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Instrument :
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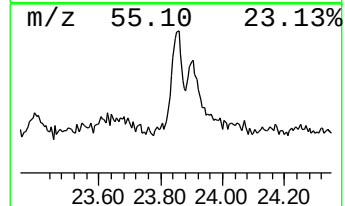
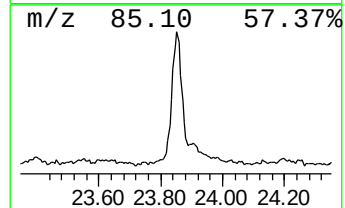
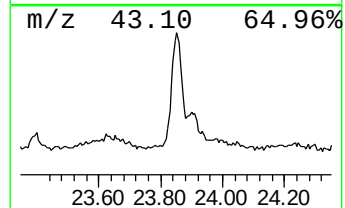
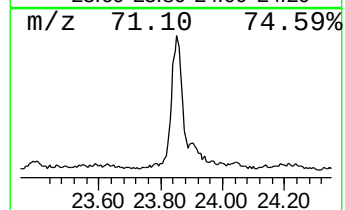
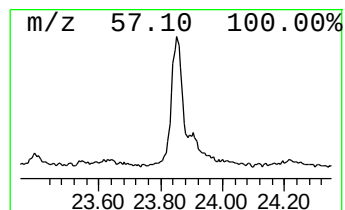
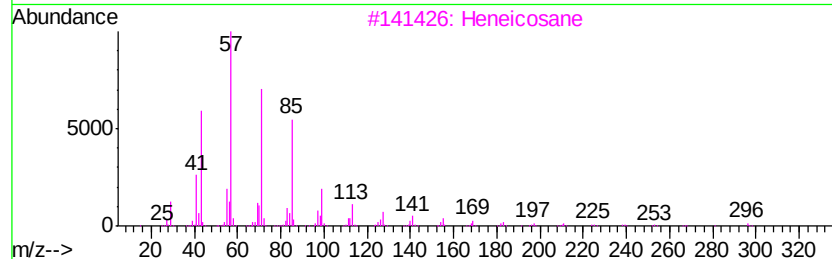
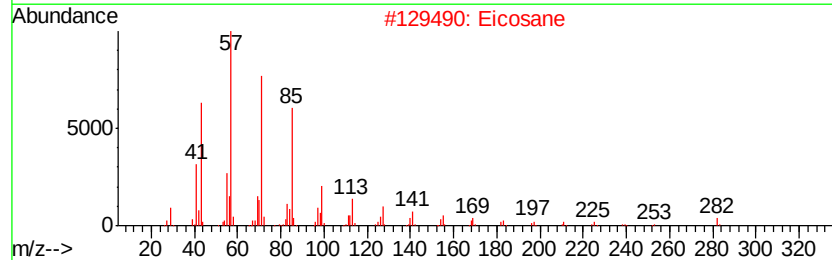
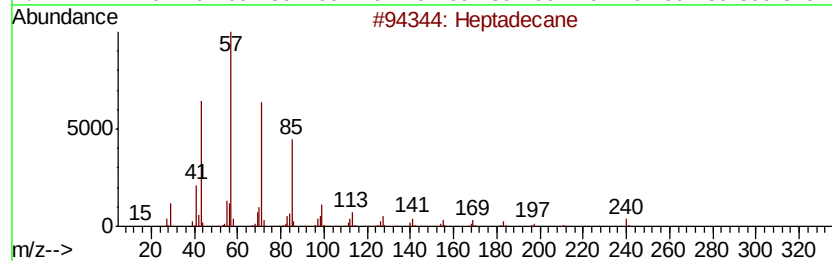
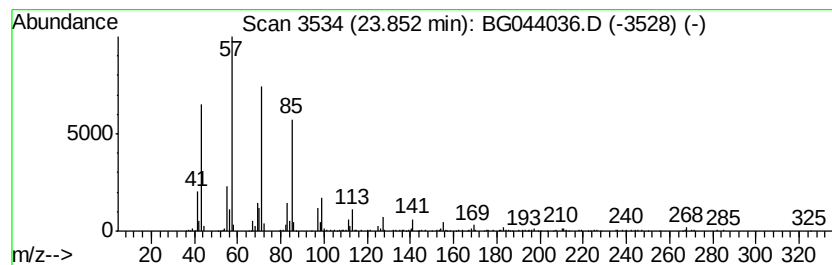
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TIC Library : C:\Database\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 14 Heptadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.85	4.94 ng	438045	Perylene-d12	24.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	95
2		Eicosane	282	C20H42	000112-95-8	94
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\BG011420\
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Misc :
ALS Vial : 5 Sample Multiplier: 1

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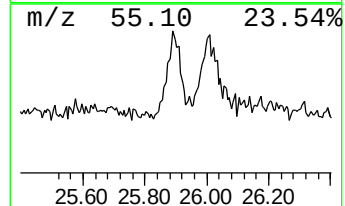
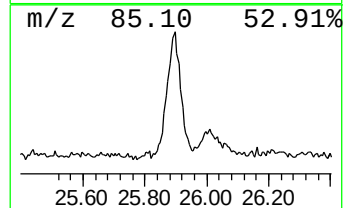
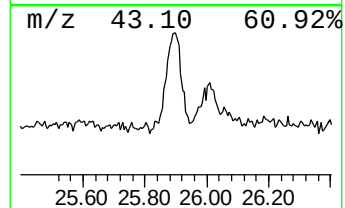
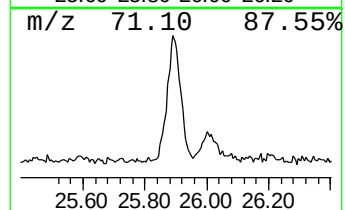
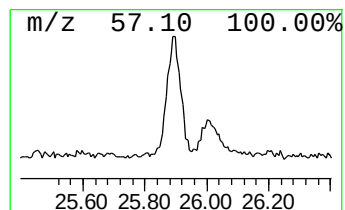
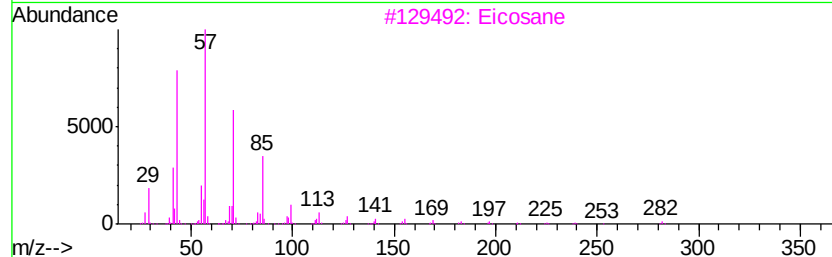
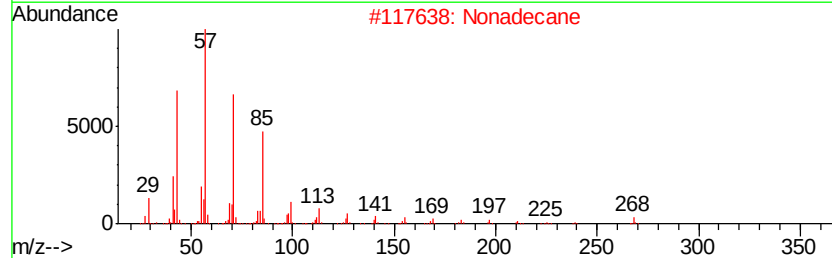
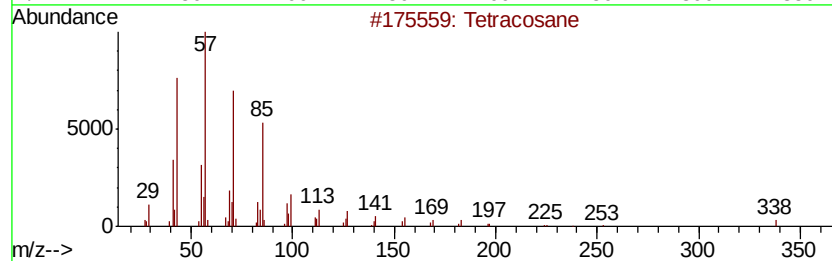
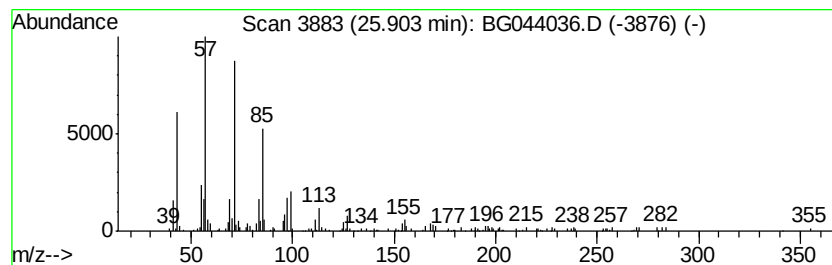
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Tetracosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.90	3.21 ng	284402	Perylene-d12	24.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	94
2		Nonadecane	268	C19H40	000629-92-5	93
3		Eicosane	282	C20H42	000112-95-8	89
4		Hexacosane	366	C26H54	000630-01-3	87
5		2-methyloctacosane	408	C29H60	1000376-72-8	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG011420\
 Data File : BG044036.D
 Acq On : 14 Jan 2020 14:49
 Operator : JU
 Sample : L1108-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 INWOOD-BH-01-A

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
Butane, 2-methoxy...	3.14	34.9	ng	1238760	1	7.95	710196	20.0
2-Pentanone, 4-hy...	5.06	15.2	ng	538343	1	7.95	710196	20.0
unknown7.48	7.48	74.0	ng	2629490	1	7.95	710196	20.0
n-Hexadecanoic acid	18.22	3.6	ng	283202	4	17.32	1578920	20.0
Heneicosane	20.66	46.1	ng	4219740	5	21.57	1831830	20.0
Heptadecyl heptaf...	21.24	8.7	ng	794117	5	21.57	1831830	20.0
Diethylene glycol...	21.30	2.3	ng	211452	5	21.57	1831830	20.0
10-Methylnonadecane	21.78	14.1	ng	1289290	5	21.57	1831830	20.0
Pentadecane, 8-he...	22.38	5.4	ng	494810	5	21.57	1831830	20.0
Eicosane	23.06	5.2	ng	477957	5	21.57	1831830	20.0
Squalene	23.24	7.1	ng	626976	6	24.70	1772710	20.0
Heptadecane	23.85	4.9	ng	438045	6	24.70	1772710	20.0
Tetracosane	25.90	3.2	ng	284402	6	24.70	1772710	20.0