

Data Path : Z:\HPCHEM1\BNA G\DATA\BG120615\
 Data File : BG019969.D
 Acq On : 6 Dec 2015 16:23
 Operator : UM/NP
 Sample : G4630-18
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-12

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:\HPCHEM1\BNA_G\METHODS\8270-BG120215.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	5.190	394	400	406	rBV	42262	64982	2.02%	0.471%
2	5.660	472	480	492	rBV	290676	469377	14.62%	3.400%
3	7.258	745	752	763	rBV	200701	338263	10.54%	2.450%
4	7.640	810	817	827	rBV	547666	933568	29.08%	6.762%
5	8.116	891	898	904	rBV	122564	207406	6.46%	1.502%
6	8.439	945	953	961	rBV	504613	859536	26.77%	6.226%
7	9.279	1087	1096	1105	rBV	440558	776968	24.20%	5.628%
8	10.930	1371	1377	1385	rVB	173060	303998	9.47%	2.202%
9	13.368	1784	1792	1800	rBV	1285344	1980534	61.69%	14.346%
10	14.737	2020	2025	2034	rVB2	281312	469743	14.63%	3.403%
11	16.224	2270	2278	2293	rBV	1078947	1641510	51.13%	11.890%
12	17.481	2485	2492	2500	rBV2	398369	607234	18.91%	4.399%
13	18.357	2636	2641	2647	rBV2	53264	74639	2.32%	0.541%
14	19.620	2852	2856	2863	rBV2	50528	77712	2.42%	0.563%
15	20.084	2929	2935	2942	rBV	2433093	3210403	100.00%	23.255%
16	20.766	3043	3051	3056	rBV	41024	71063	2.21%	0.515%
17	20.977	3084	3087	3092	rBV2	27651	40175	1.25%	0.291%
18	21.030	3092	3096	3100	rVV	42598	59647	1.86%	0.432%
19	21.388	3154	3157	3165	rVB8	20204	37468	1.17%	0.271%
20	21.535	3174	3182	3190	rBV	37903	81058	2.52%	0.587%
21	21.753	3213	3219	3228	rVB	478999	782828	24.38%	5.670%
22	25.031	3766	3777	3791	rVB3	252746	717270	22.34%	5.196%

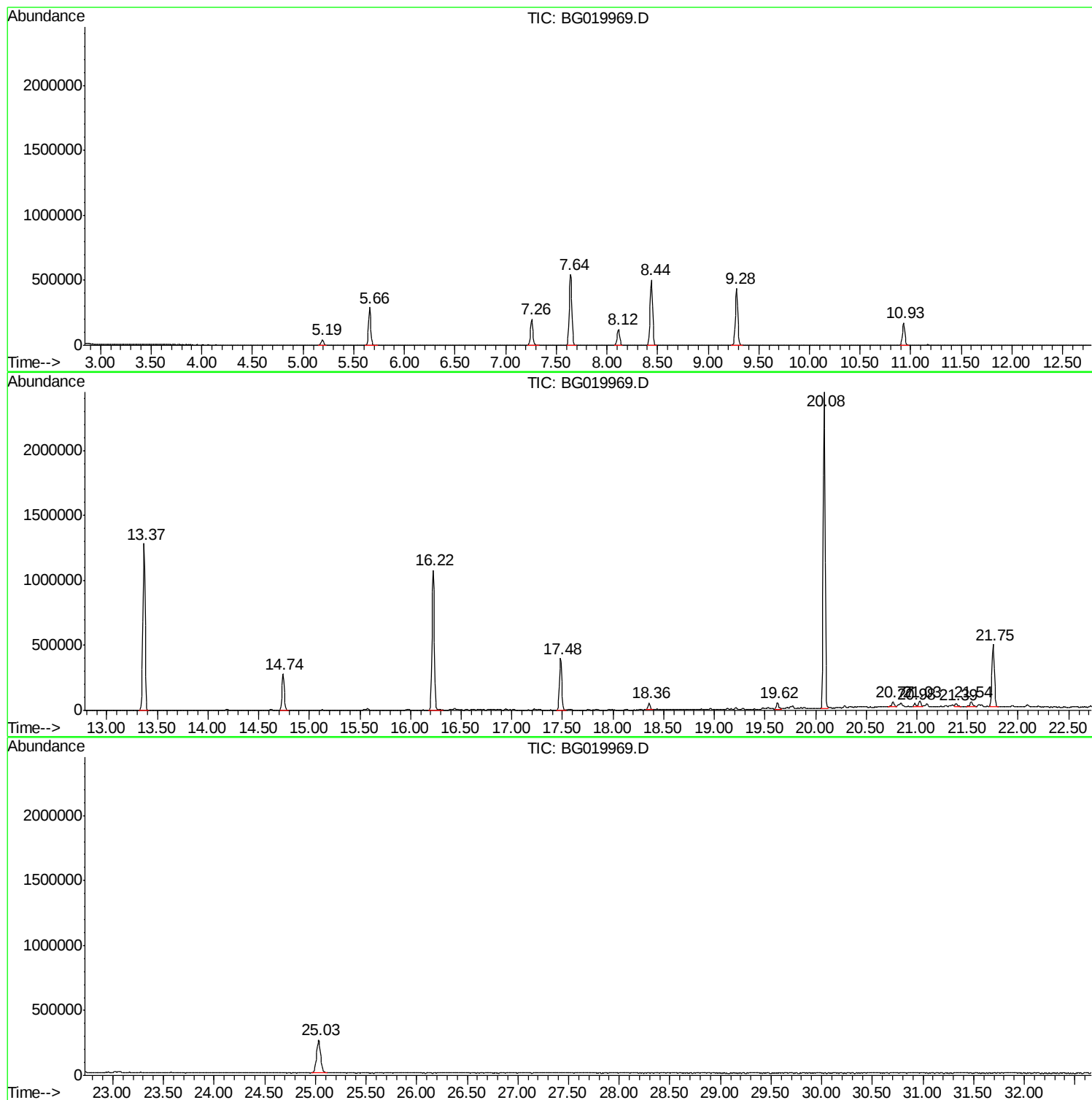
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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Instrument :
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ClientSampled :
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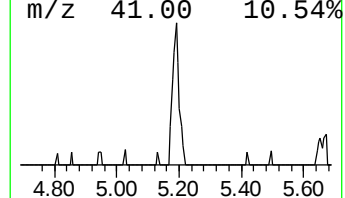
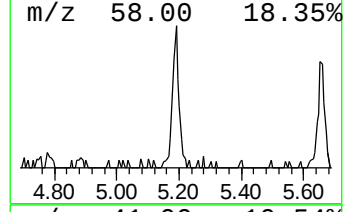
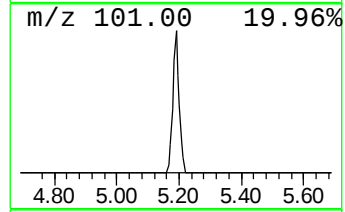
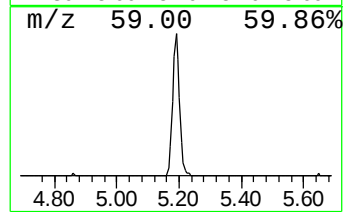
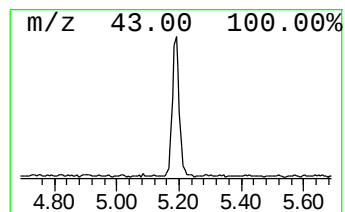
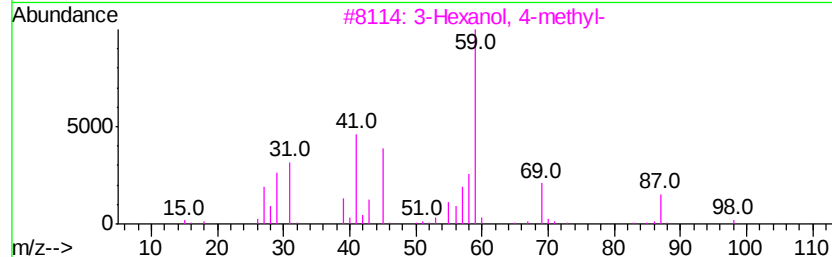
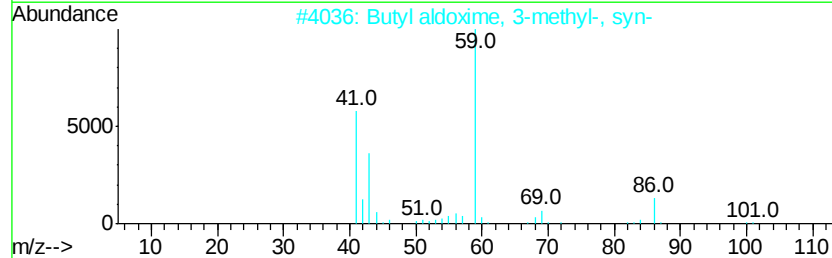
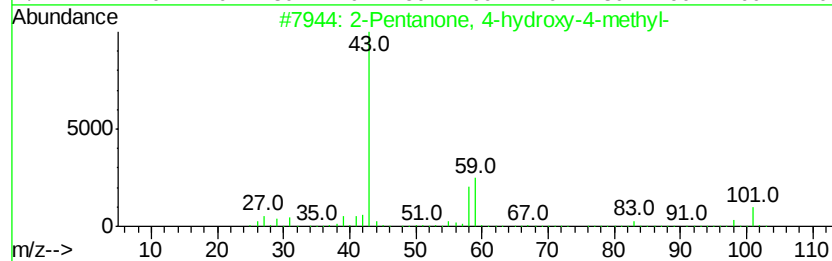
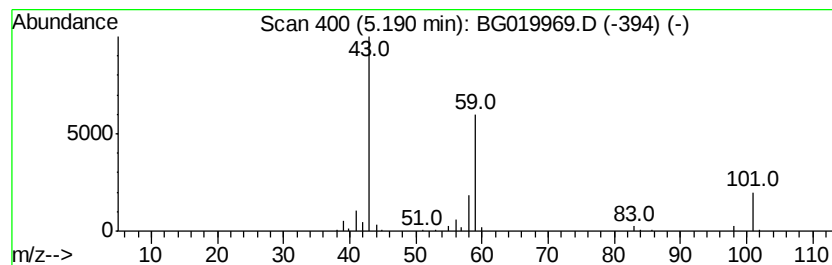
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	6.27 ng	64982	1,4-Dichlorobenzene-d4	8.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	35
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
5		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	17



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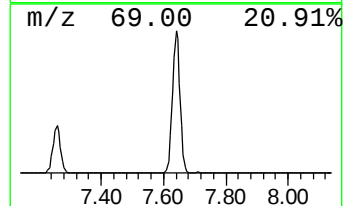
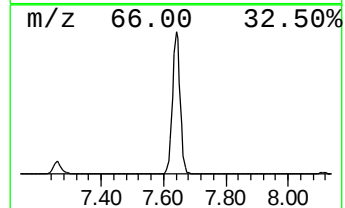
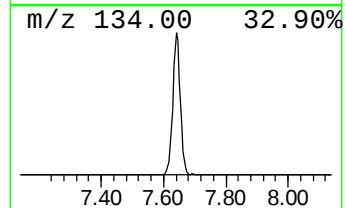
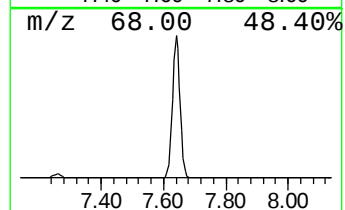
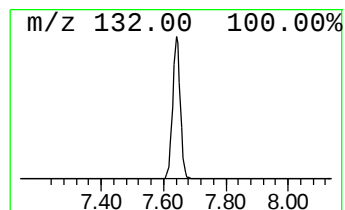
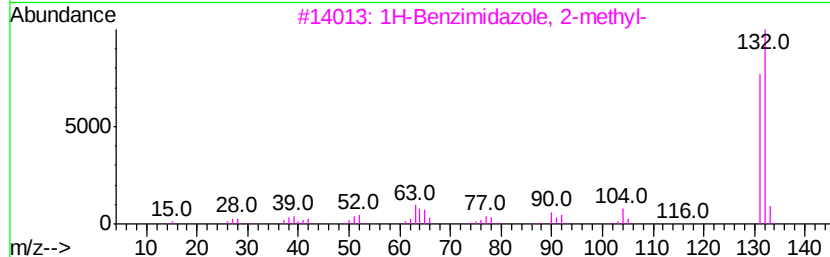
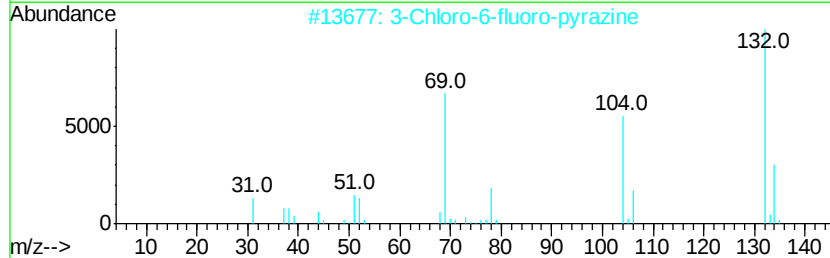
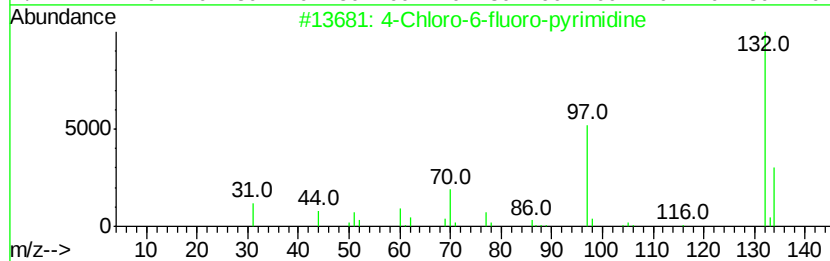
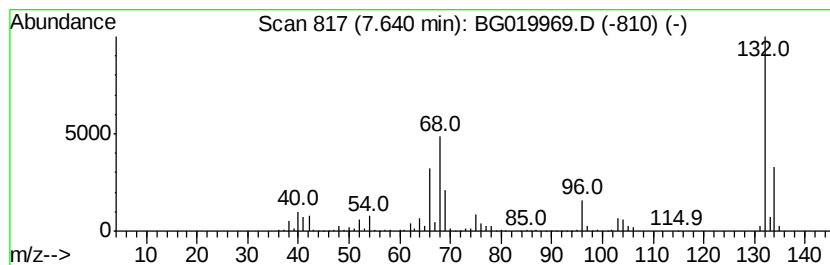
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TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown7.64 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.64	90.02 ng	933568	1,4-Dichlorobenzene-d4	8.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	14
5		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	13



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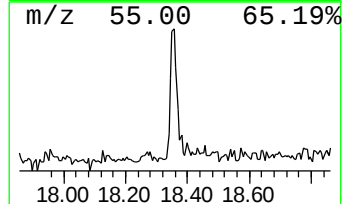
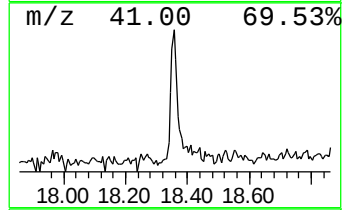
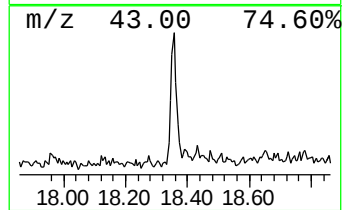
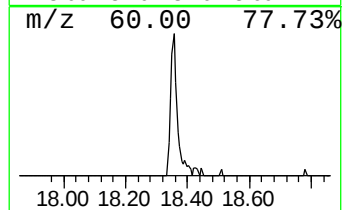
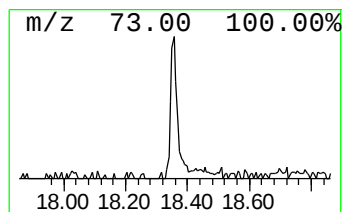
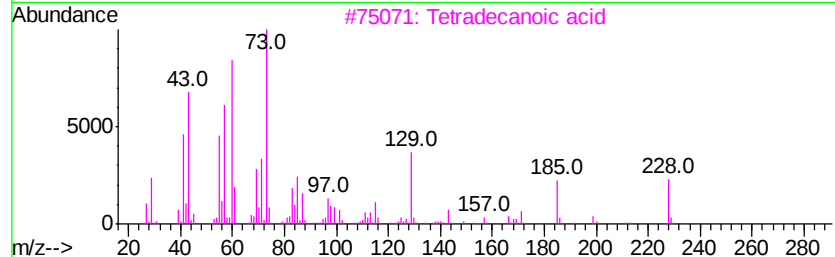
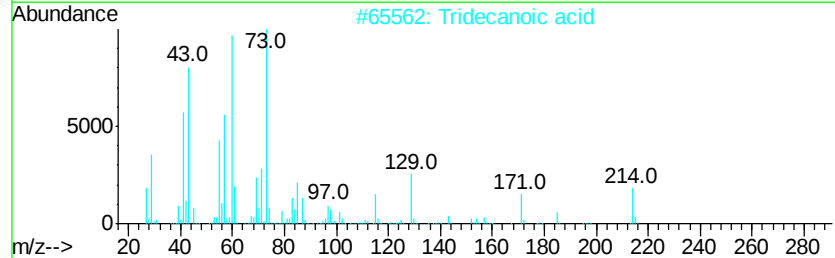
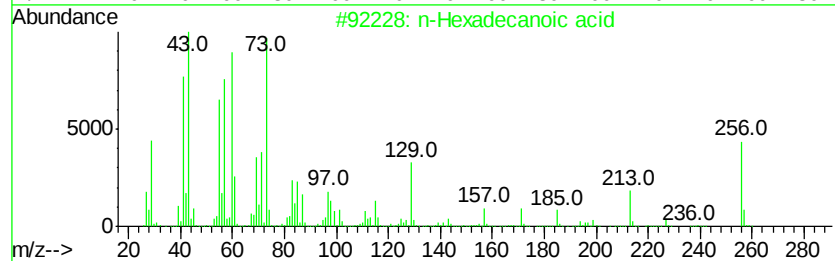
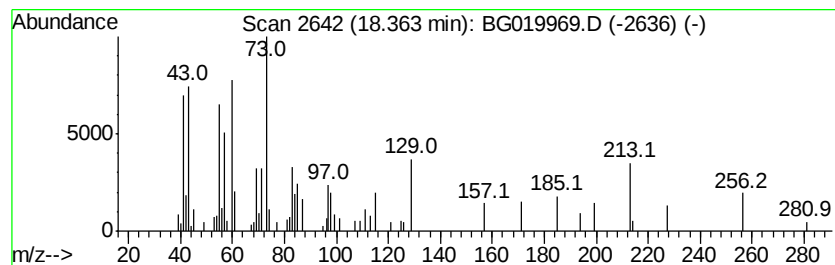
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.36	2.46 ng	74639	Phenanthrene-d10	17.48

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2		Tridecanoic acid	214	C13H26O2	000638-53-9	81
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	70
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	50
5		Undecanoic acid	186	C11H22O2	000112-37-8	50



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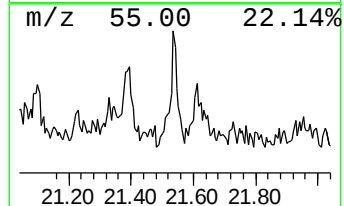
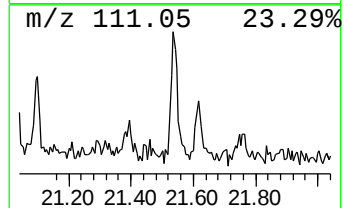
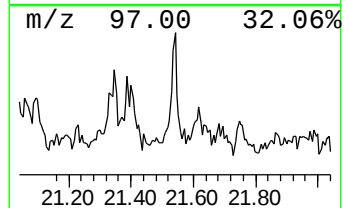
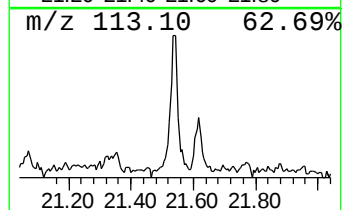
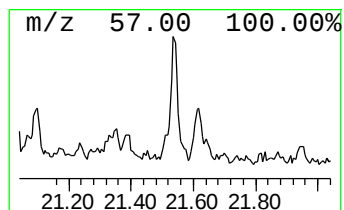
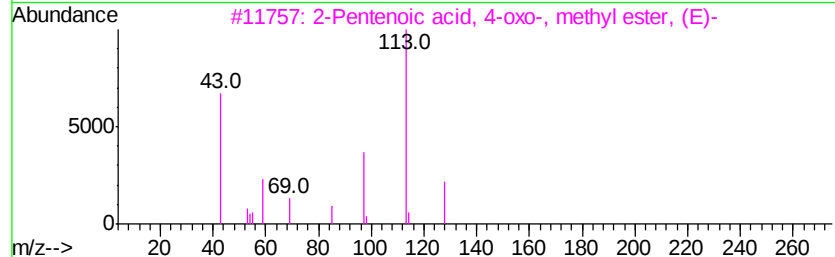
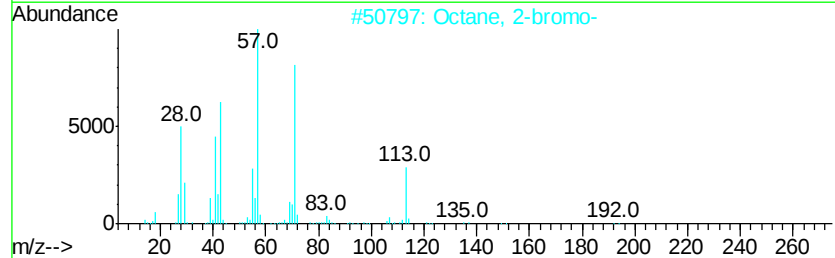
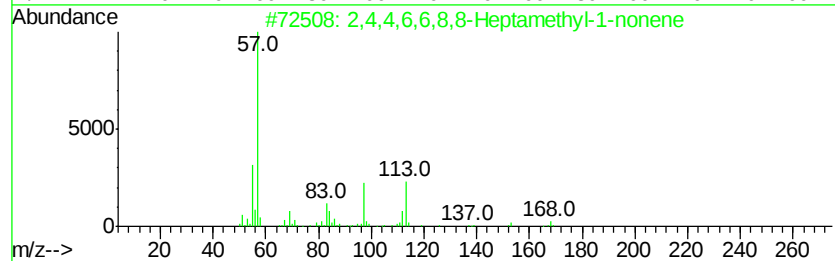
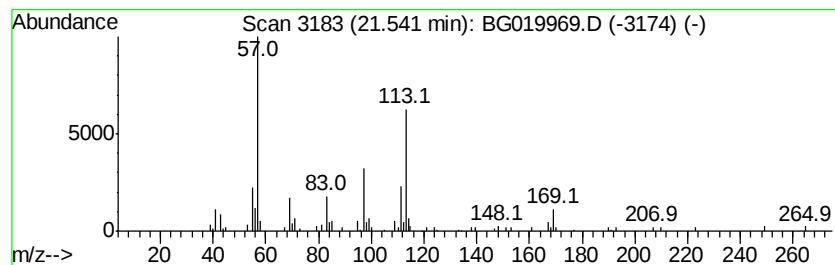
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Peak Number 5 unknown21.54 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.54	2.07 ng	81058	Chrysene-d12	21.75

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,4,6,6,8,8-Heptamethyl-1-nonene	224	C16H32	015796-04-0	43		
2	Octane, 2-bromo-	192	C8H17Br	000557-35-7	43		
3	2-Pentenoic acid, 4-oxo-, methyl...	128	C6H8O3	002833-24-1	43		
4	Silane, butyldichloromethyl-	170	C5H12Cl2Si	018147-23-4	27		
5	Pyrazole, 3-nitro-	113	C3H3N3O2	026621-44-3	27		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.19	6.3	ng	64982	1	8.12	207406	20.0
unknown7.64	7.64	90.0	ng	933568	1	8.12	207406	20.0
n-Hexadecanoic acid	18.36	2.5	ng	74639	4	17.48	607234	20.0
unknown21.54	21.54	2.1	ng	81058	5	21.75	782828	20.0