

Data Path : Z:\HPCHEM1\BNA G\DATA\BG021816\
 Data File : BG020958.D
 Acq On : 19 Feb 2016 5:38
 Operator : SJ/UM
 Sample : H1488-01
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 P001-SS016-5460-01

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:\HPCHEM1\BNA_G\METHODS\8270-BG021116.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	3.090	37	43	63	rVB	1081275	1515689	98.61%	14.660%
2	3.237	63	68	77	rBV	23962	41449	2.70%	0.401%
3	5.311	415	421	435	rVB	78765	122164	7.95%	1.182%
4	5.793	496	503	512	rBV	382092	599677	39.01%	5.800%
5	7.402	769	777	790	rBV	401453	693131	45.09%	6.704%
6	7.784	835	842	855	rVB	426405	727391	47.32%	7.036%
7	8.260	916	923	929	rVB	73537	123920	8.06%	1.199%
8	8.583	970	978	987	rVB	314767	548363	35.68%	5.304%
9	9.429	1114	1122	1130	rBV	249450	440355	28.65%	4.259%
10	11.086	1395	1404	1411	rBV	126072	224322	14.59%	2.170%
11	13.500	1807	1815	1824	rBV	899511	1330659	86.57%	12.871%
12	14.311	1947	1953	1960	rVB	76365	105700	6.88%	1.022%
13	14.869	2042	2048	2056	rVB2	207356	325581	21.18%	3.149%
14	15.685	2183	2187	2192	rVB	20537	24547	1.60%	0.237%
15	16.349	2294	2300	2312	rVB	541856	781675	50.86%	7.561%
16	16.543	2328	2333	2336	rBV	19042	27029	1.76%	0.261%
17	17.606	2508	2514	2520	rBV	235046	337766	21.97%	3.267%
18	20.185	2947	2953	2959	rBV	1236582	1537063	100.00%	14.867%
19	21.483	3169	3174	3186	rBV	64787	101243	6.59%	0.979%
20	21.830	3227	3233	3236	rBV5	9103	15384	1.00%	0.149%
21	21.883	3236	3242	3248	rVB	217482	359085	23.36%	3.473%
22	25.249	3805	3815	3829	rVB3	122205	356479	23.19%	3.448%

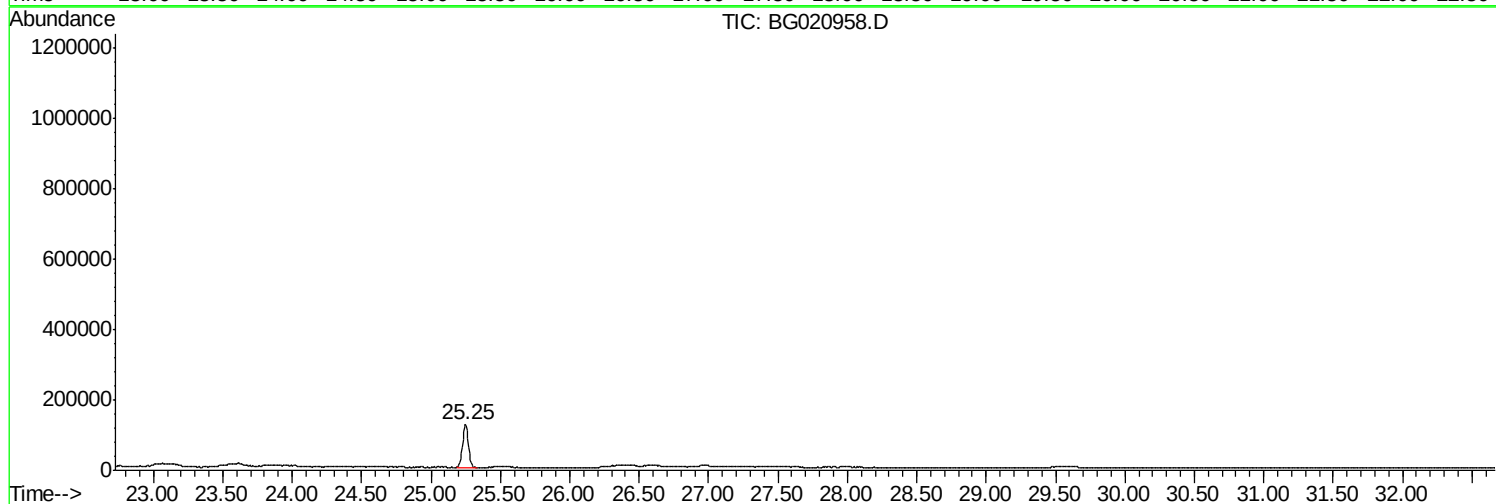
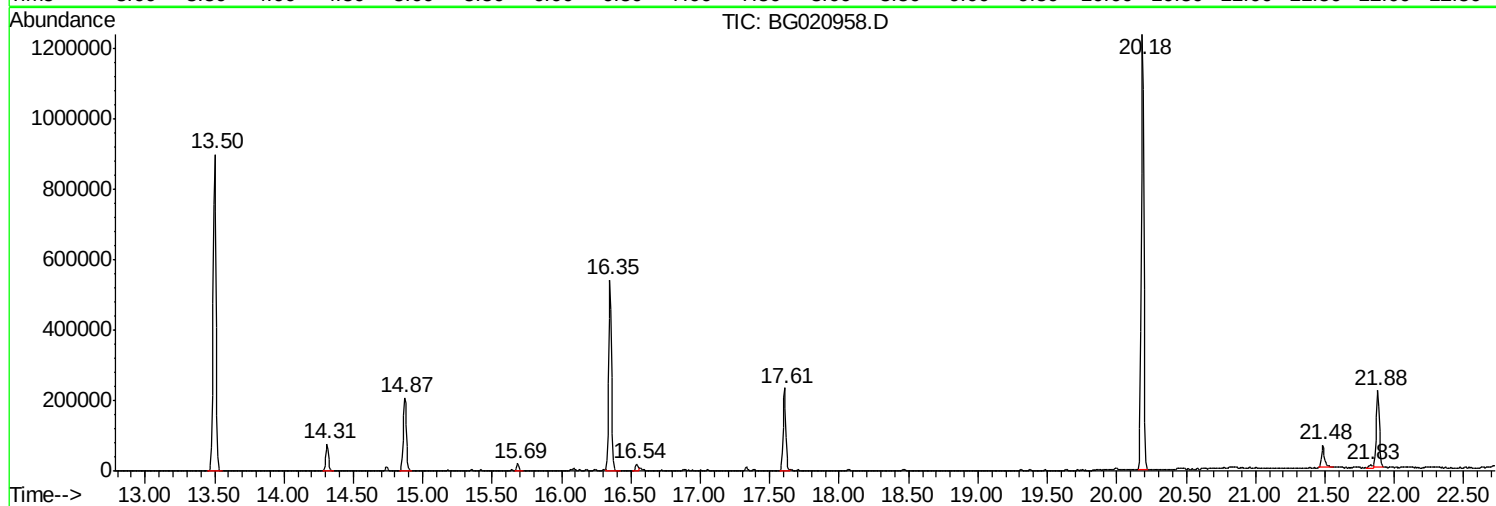
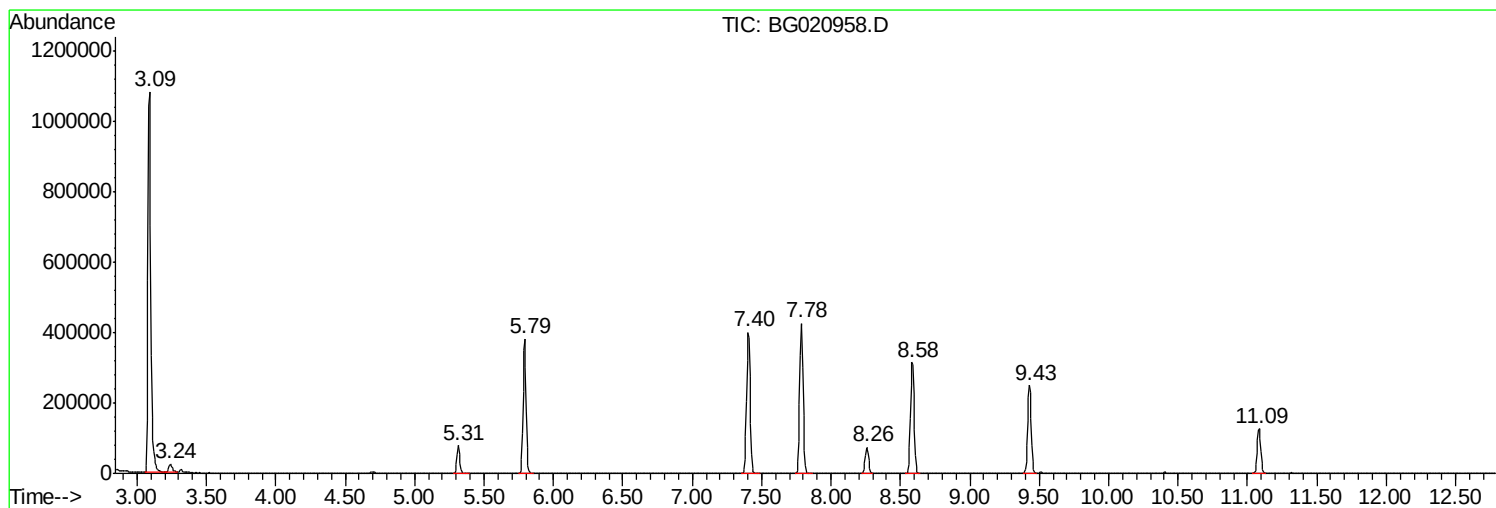
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG021116.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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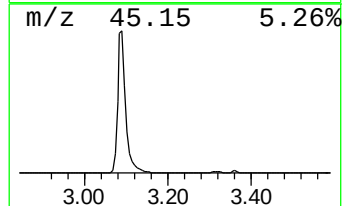
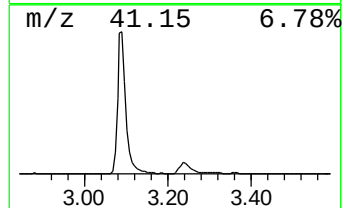
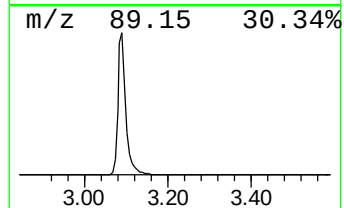
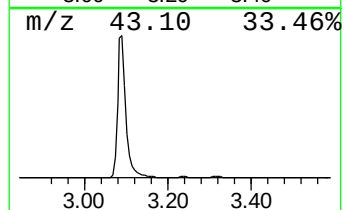
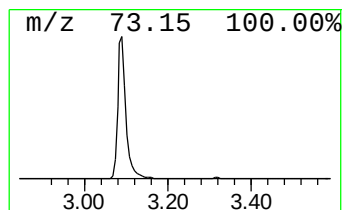
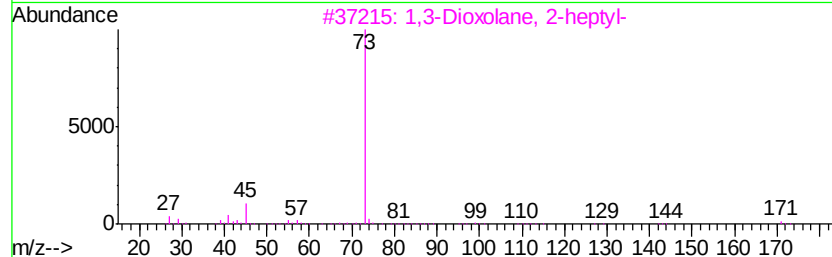
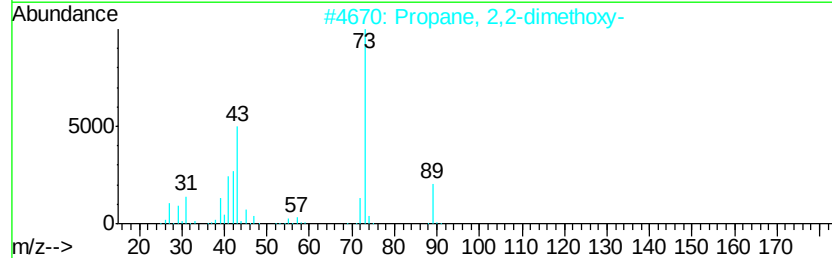
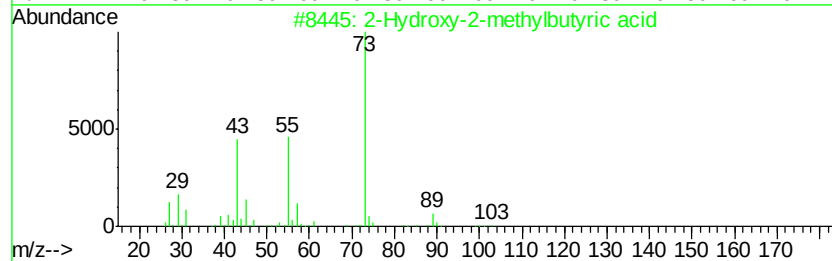
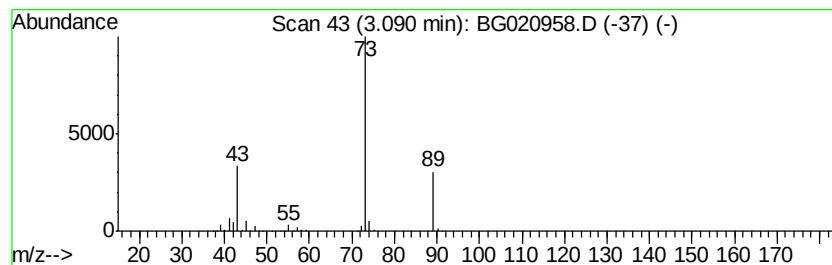
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG021116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown3.09 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.09	244.62 ng	1515690	1,4-Dichlorobenzene-d4	8.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	28
2		Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	9
3		1,3-Dioxolane, 2-heptyl-	172	C10H20O2	004359-57-3	9
4		N-Ethylformamide	73	C3H7NO	000627-45-2	7
5		Methane, isothiocyanato-	73	C2H3NS	000556-61-6	7



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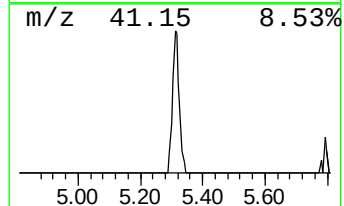
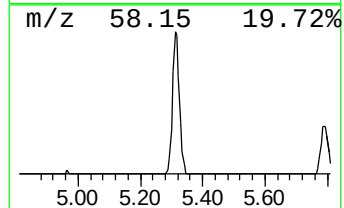
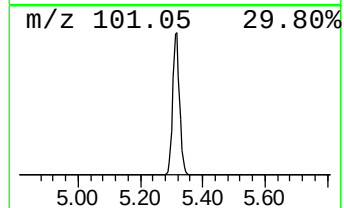
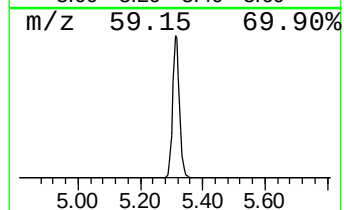
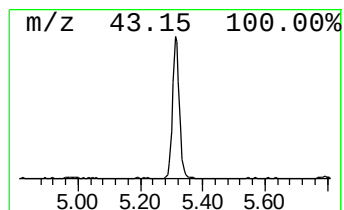
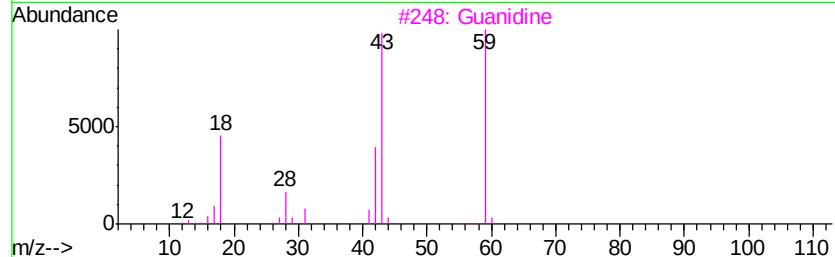
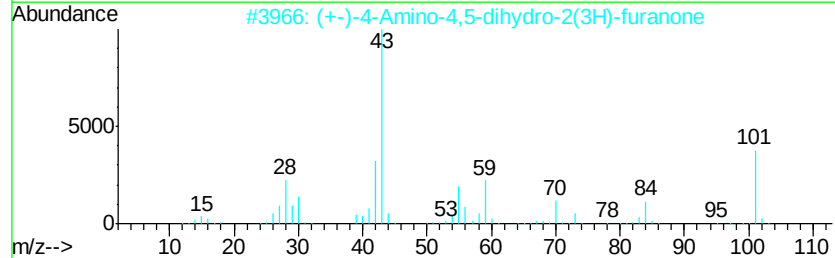
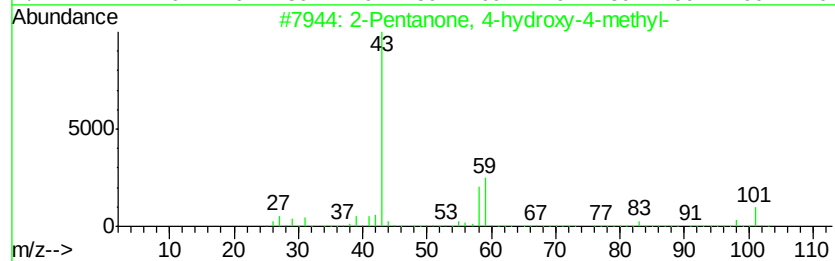
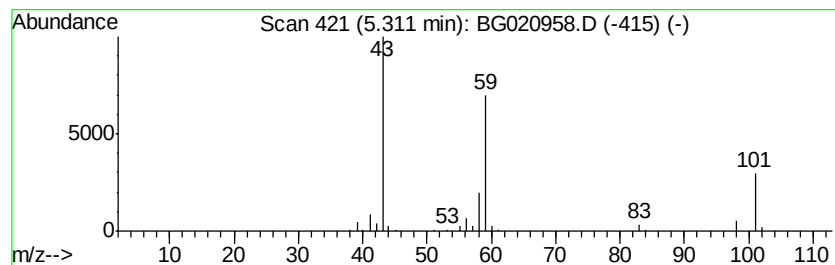
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.31	19.72 ng	122164	1,4-Dichlorobenzene-d4	8.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2		(+)-4-Amino-4,5-dihydro-2(3H)-furanone	101	C4H7NO2	016504-58-8	43
3		Guanidine	59	CH5N3	000113-00-8	38
4		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	12
5		5-Hexen-2-one	98	C6H10O	000109-49-9	9



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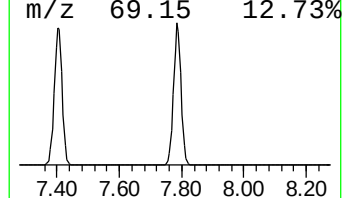
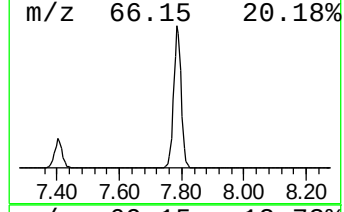
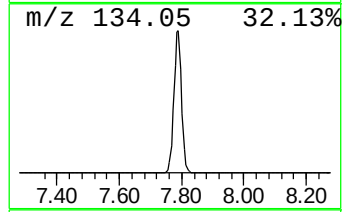
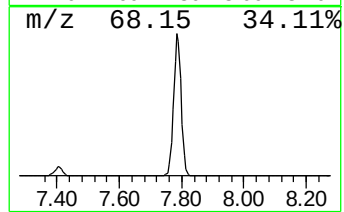
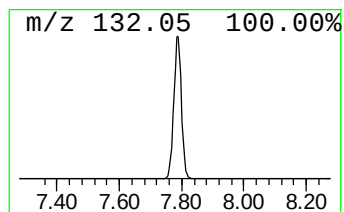
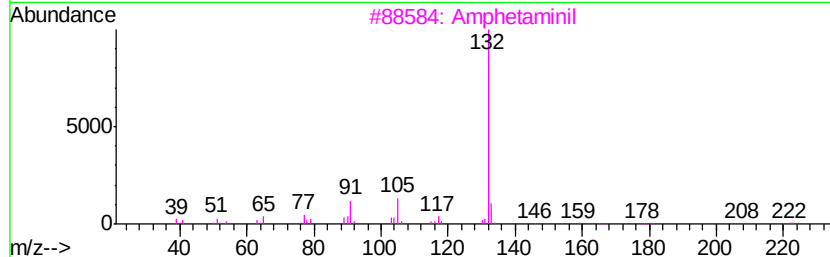
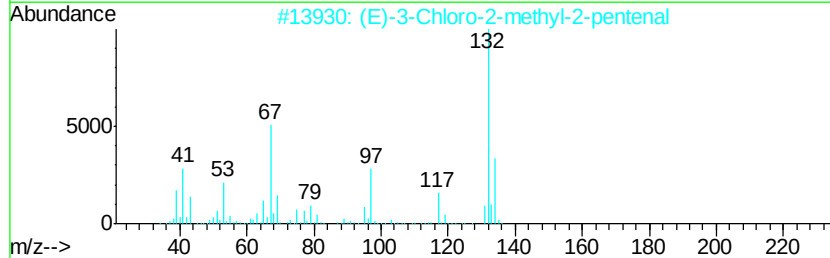
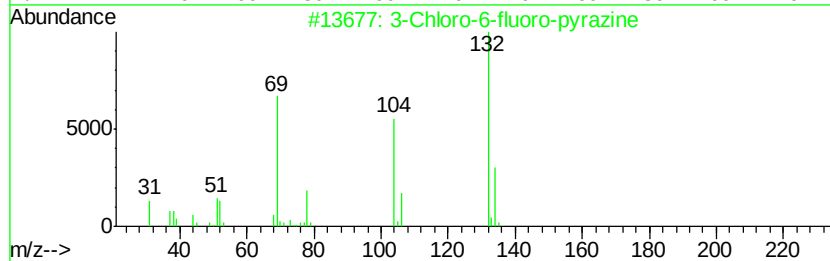
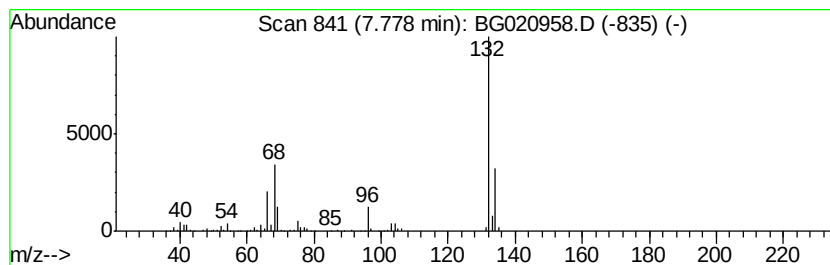
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Peak Number 4 unknown7.78 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.78	117.40 ng	727391	1,4-Dichlorobenzene-d4	8.25

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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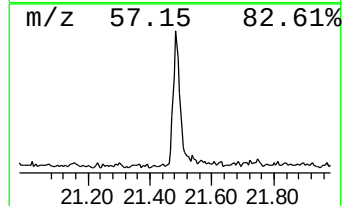
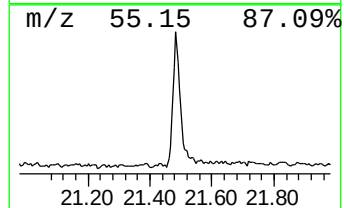
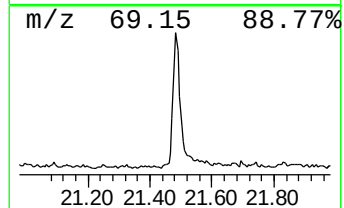
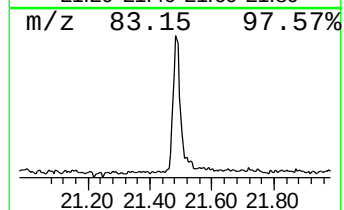
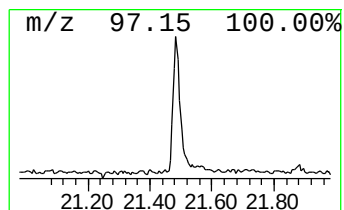
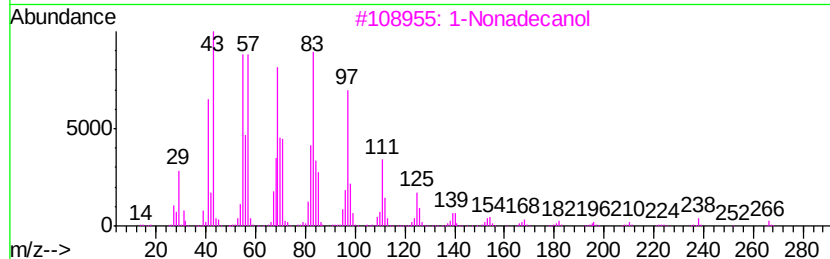
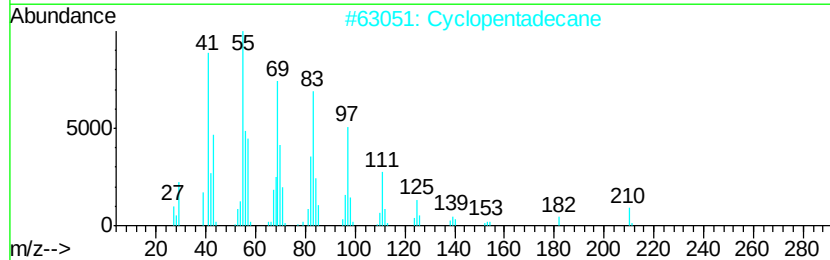
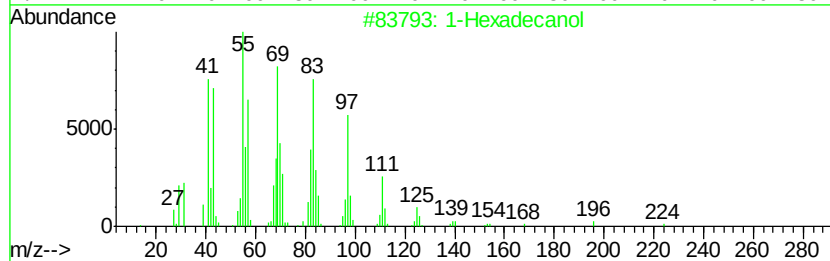
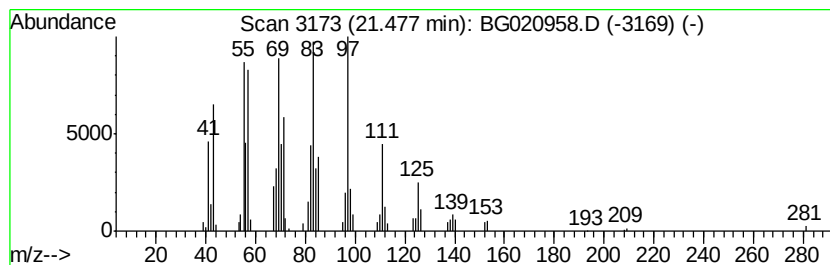
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Peak Number 6 1-Hexadecanol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.48	5.64 ng	101243	Chrysene-d12	21.88	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexadecanol	242	C16H34O	036653-82-4	91
2	Cyclopentadecane	210	C15H30	000295-48-7	91
3	1-Nonadecanol	284	C19H40O	001454-84-8	91
4	Cyclohexadecane	224	C16H32	000295-65-8	91
5	2- Chloropropionic acid, hexadec...	332	C19H37ClO2	086711-81-1	90



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown3.09	3.09	244.6	ng	1515690	1	8.25	123920	20.0
2-Pentanone, 4-hy...	5.31	19.7	ng	122164	1	8.25	123920	20.0
unknown7.78	7.78	117.4	ng	727391	1	8.25	123920	20.0
1-Hexadecanol	21.48	5.6	ng	101243	5	21.88	359085	20.0