

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_G\DATA\BG011518\
 Data File : BG032046.D
 Acq On : 15 Jan 2018 23:51
 Operator : SJ/JU
 Sample : J1118-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DUP-2

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-BG011218.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	3.406	14	20	30	rBV3	41012	89590	2.42%	0.503%
2	3.494	31	35	46	rVB	35171	56743	1.53%	0.318%
3	5.686	402	408	415	rVB	35833	57008	1.54%	0.320%
4	6.191	487	494	506	rBV	636452	1112892	30.01%	6.245%
5	7.871	772	780	792	rBV	634739	1207186	32.55%	6.774%
6	8.277	841	849	863	rBV	685781	1246473	33.61%	6.994%
7	8.764	926	932	941	rBV	137136	249664	6.73%	1.401%
8	9.105	982	990	1002	rVB	512032	974091	26.27%	5.466%
9	9.963	1128	1136	1147	rBV	383999	709058	19.12%	3.979%
10	11.649	1415	1423	1432	rBV	187231	353105	9.52%	1.981%
11	14.029	1821	1828	1840	rVB	1283551	2080886	56.12%	11.677%
12	14.822	1955	1963	1971	rBV	101272	154365	4.16%	0.866%
13	15.398	2054	2061	2071	rVB2	335227	573952	15.48%	3.221%
14	16.873	2305	2312	2324	rBV	1399886	2135291	57.58%	11.982%
15	18.130	2520	2526	2537	rBV2	522791	852954	23.00%	4.786%
16	20.662	2951	2957	2969	rBV	2676120	3708172	100.00%	20.808%
17	22.008	3180	3186	3190	rBV5	24645	44999	1.21%	0.253%
18	22.466	3257	3264	3278	rBV	587455	1114848	30.06%	6.256%
19	26.185	3885	3897	3910	rBV3	336509	1099571	29.65%	6.170%

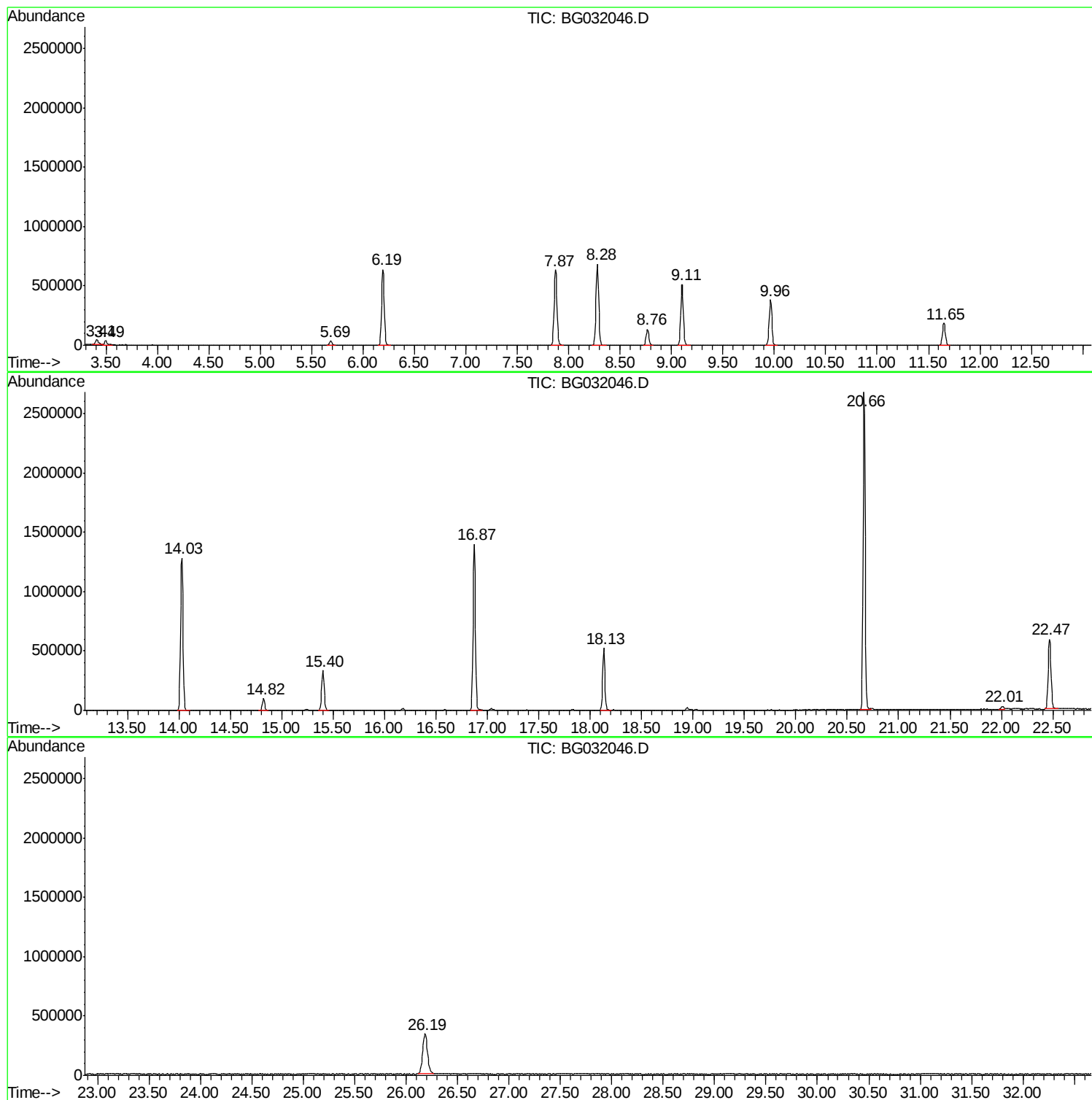
Sum of corrected areas: 17820848

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG011218.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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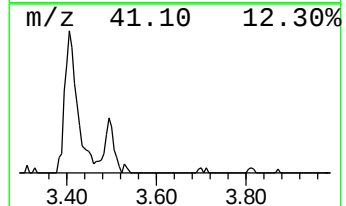
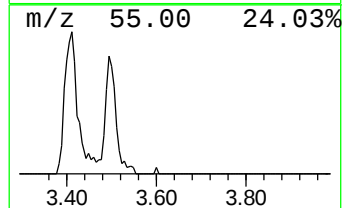
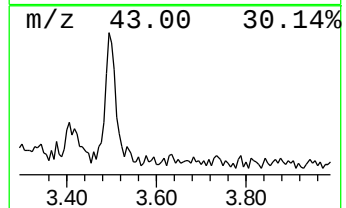
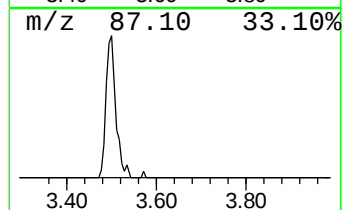
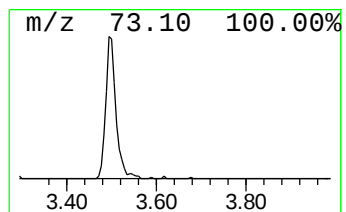
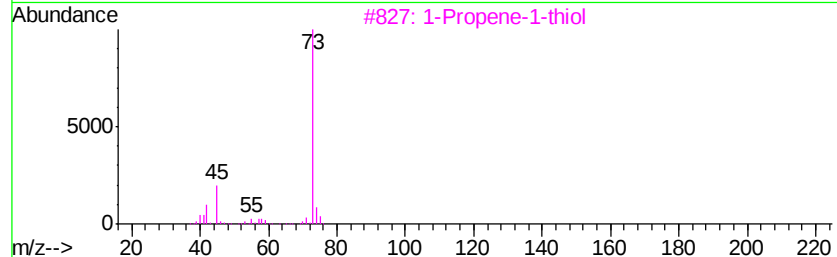
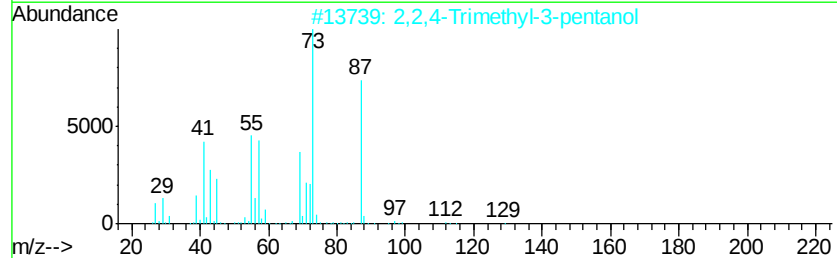
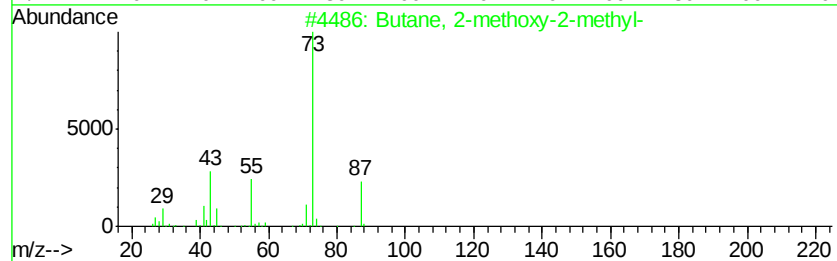
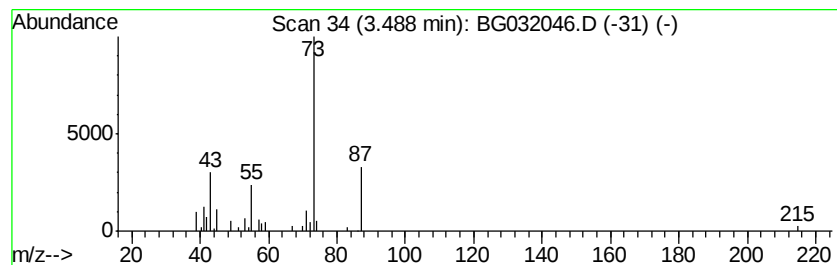
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG011218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.49	4.55 ng	56743	1,4-Dichlorobenzene-d4	8.77

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	40
3		1-Propene-1-thiol	74	C3H6S	000925-89-3	17
4		N-Ethylformamide	73	C3H7NO	000627-45-2	9
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9



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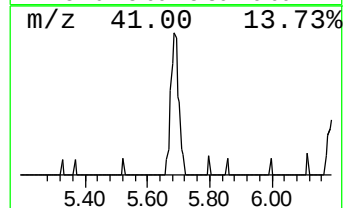
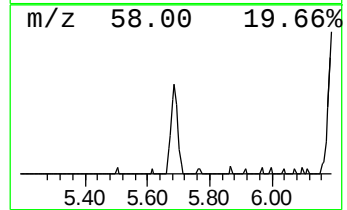
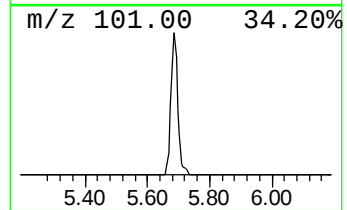
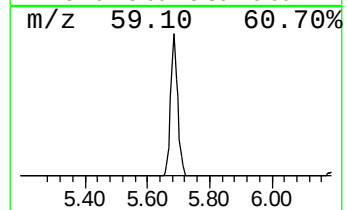
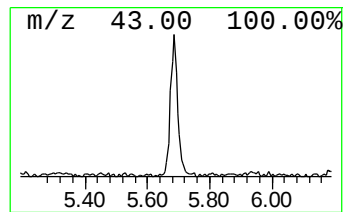
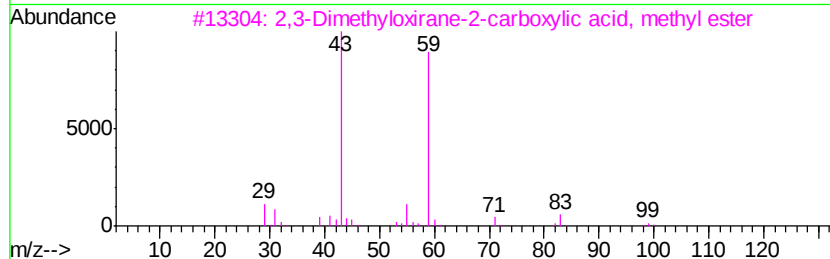
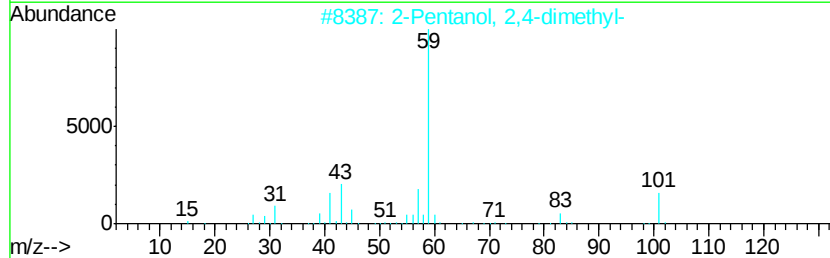
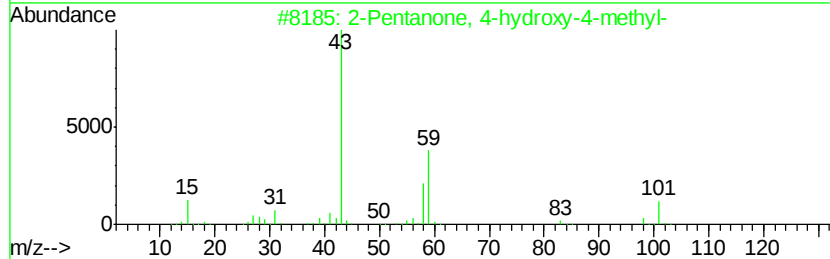
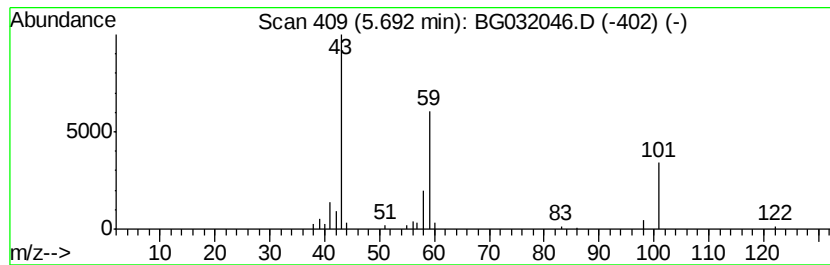
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Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.69	4.57 ng	57008	1,4-Dichlorobenzene-d4	8.77

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	78
2			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	53
3			2,3-Dimethyloxirane-2-carboxylic...	130	C6H10O3	1000306-64-4	33
4			Acetone	58	C3H6O	000067-64-1	12
5			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9



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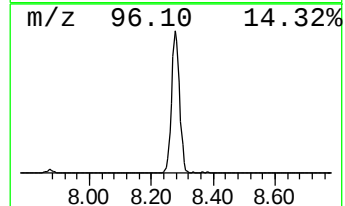
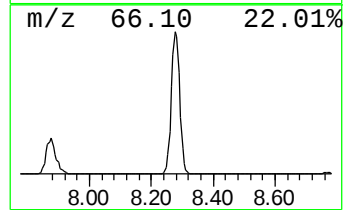
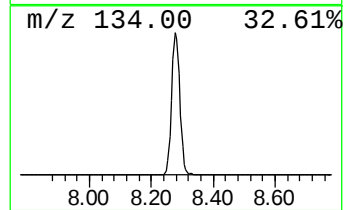
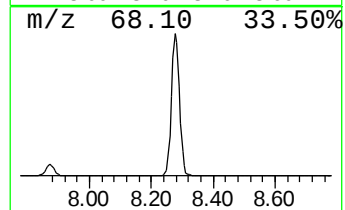
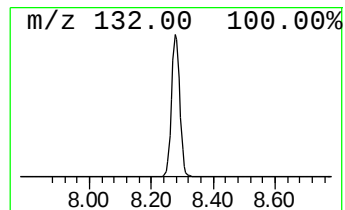
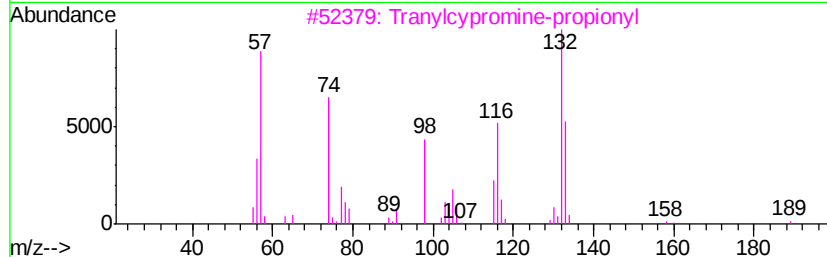
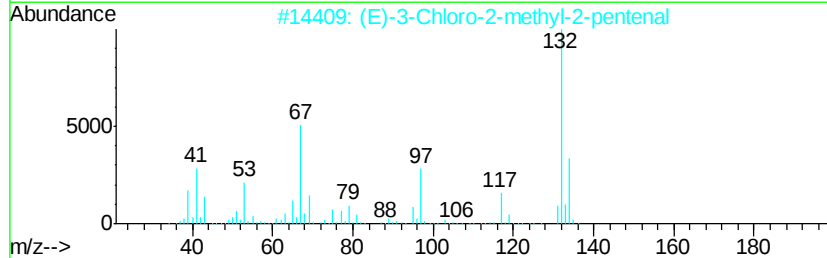
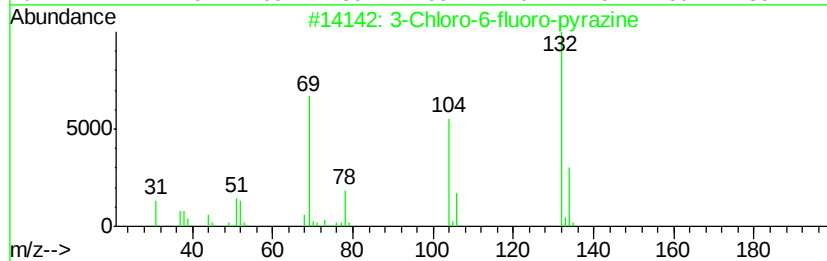
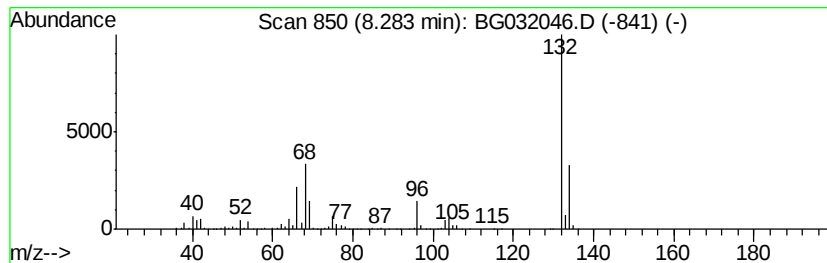
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Peak Number 4 unknown8.28 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.28	99.85 ng	1246470	1,4-Dichlorobenzene-d4	8.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.49	4.5	ng	56743	1	8.77	249664	20.0
2-Pentanone, 4-hy...	5.69	4.6	ng	57008	1	8.77	249664	20.0
unknown8.28	8.28	99.8	ng	1246470	1	8.77	249664	20.0