

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060715\
 Data File : BF079800.D
 Acq On : 7 Jun 2015 8:10
 Operator : TP/IZ
 Sample : G2508-09
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ABSLOPE-1(0-2)

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_F\METHODS\8270-BF060515.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	2.227	90	91	105	rVB	118528	307418	4.59%	1.253%
2	2.421	105	108	111	rBV	4293090	6696626	100.00%	27.302%
3	2.879	145	148	155	rVB	227290	335412	5.01%	1.367%
4	6.090	427	429	432	rBV	2094546	1812758	27.07%	7.390%
5	7.073	512	515	517	rBV	1805872	1908751	28.50%	7.782%
6	7.245	527	530	532	rBV	2099138	1963824	29.33%	8.006%
7	7.473	548	550	552	rBV	387117	343270	5.13%	1.399%
8	7.633	562	564	566	rBV	1695352	1472944	22.00%	6.005%
9	8.033	596	599	601	rBV	1064498	1054709	15.75%	4.300%
10	8.765	661	663	666	rBV	502022	450451	6.73%	1.836%
11	9.839	755	757	760	rVB	2548833	2209122	32.99%	9.006%
12	10.548	816	819	821	rVB	398637	506105	7.56%	2.063%
13	11.337	886	888	890	rBV	1463876	1250195	18.67%	5.097%
14	12.057	949	951	953	rBV	621596	565901	8.45%	2.307%
15	13.760	1097	1100	1102	rBV	2250094	2550957	38.09%	10.400%
16	15.051	1210	1213	1215	rBV	434269	570354	8.52%	2.325%
17	17.006	1380	1384	1391	rBV	300557	529543	7.91%	2.159%

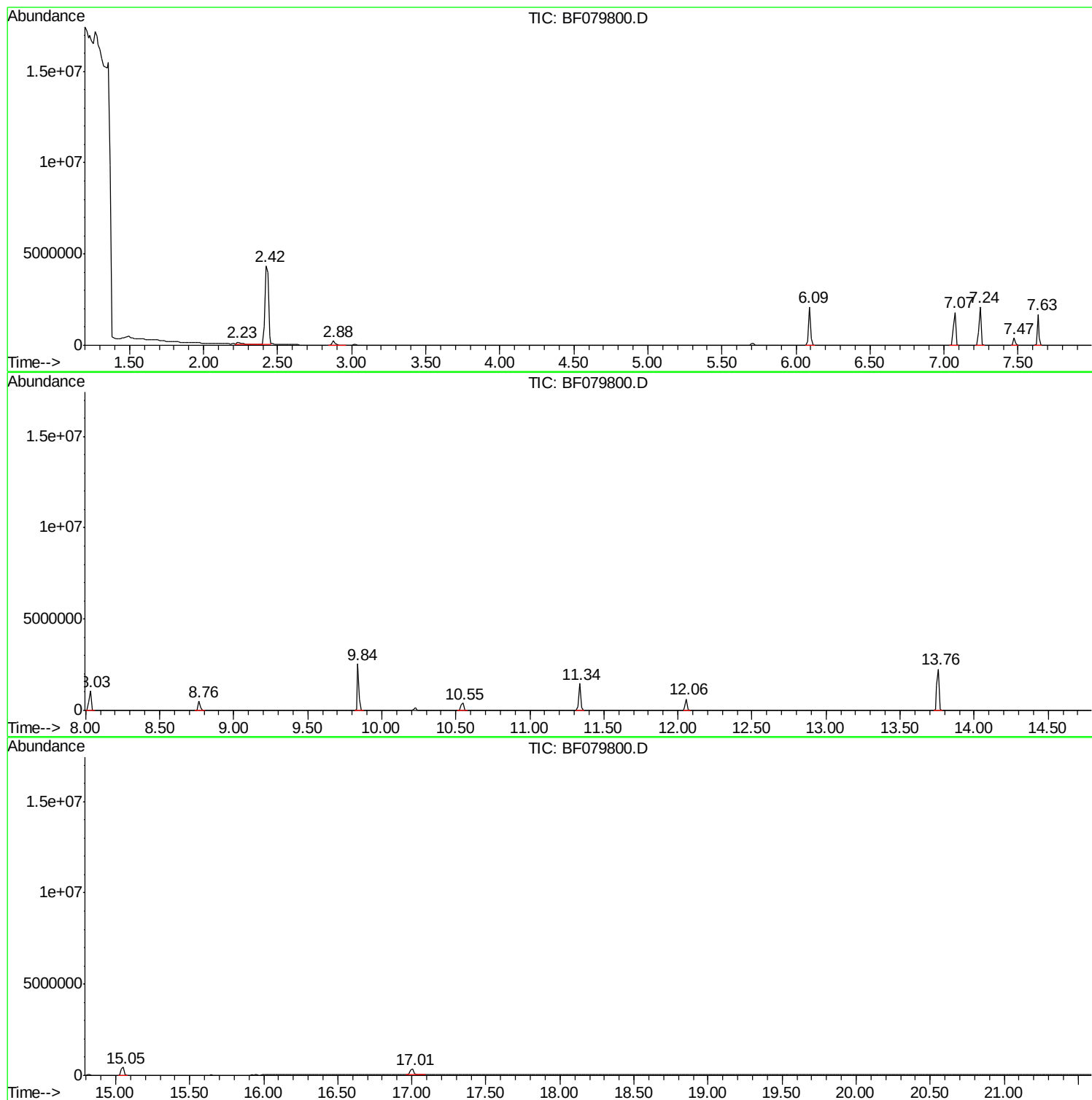
Sum of corrected areas: 24528340

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ABSLOPE-1(0-2)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060515.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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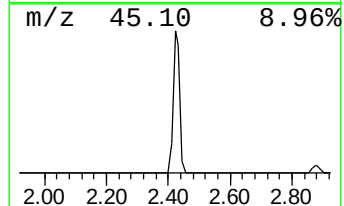
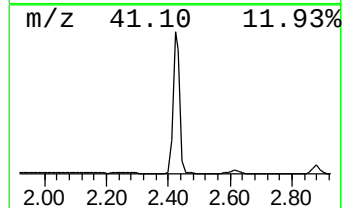
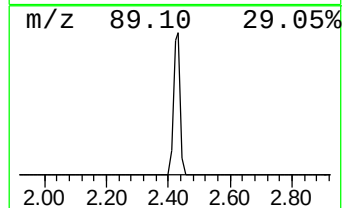
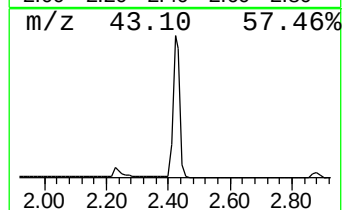
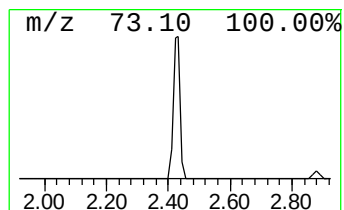
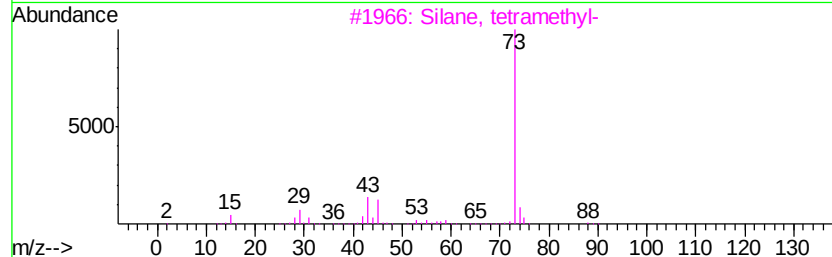
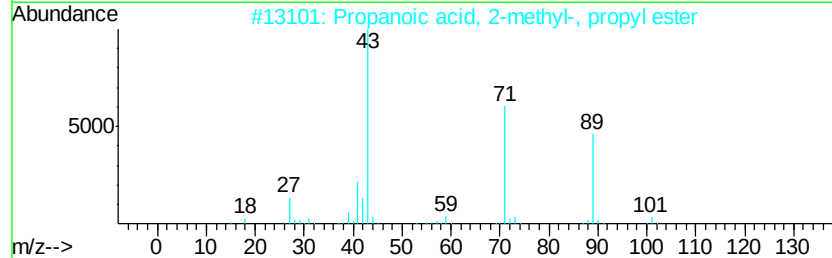
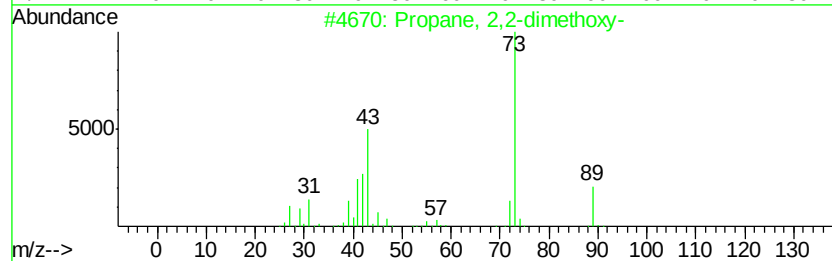
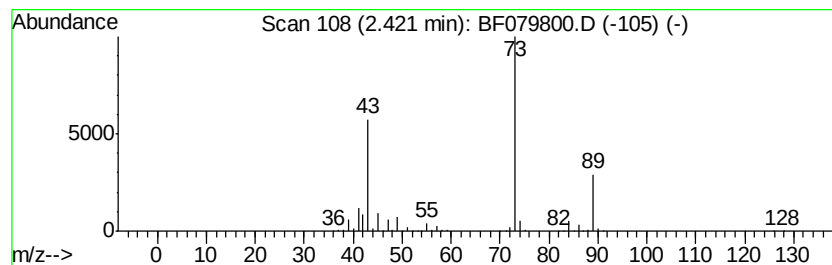
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Propane, 2,2-dimethoxy- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.42	390.17 ng	6696630	1,4-Dichlorobenzene-d4	7.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	50
2		Propanoic acid, 2-methyl-, propyl ester	130	C7H14O2	000644-49-5	12
3		Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
4		1,3-Dioxolane, 2-pentadecyl-	284	C18H36O2	004360-57-0	9
5		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9



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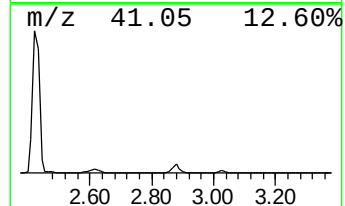
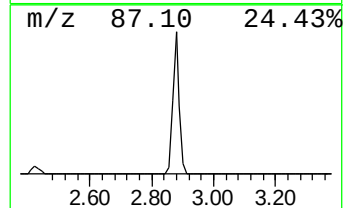
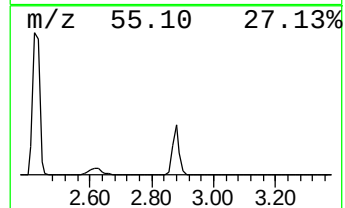
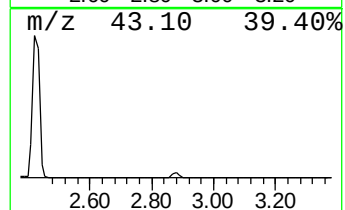
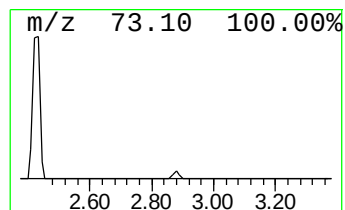
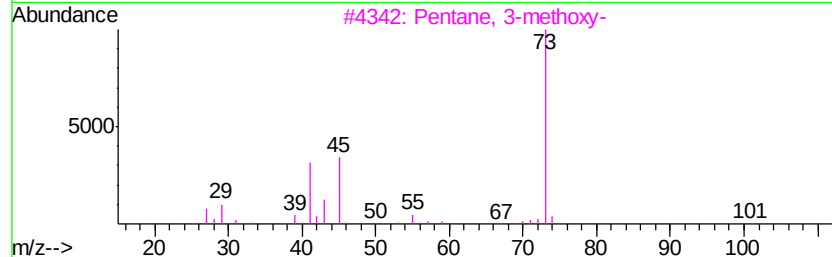
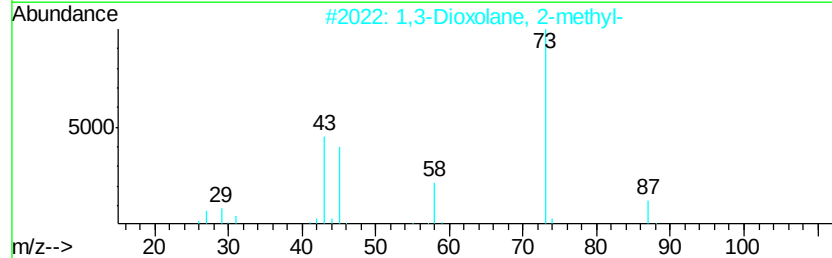
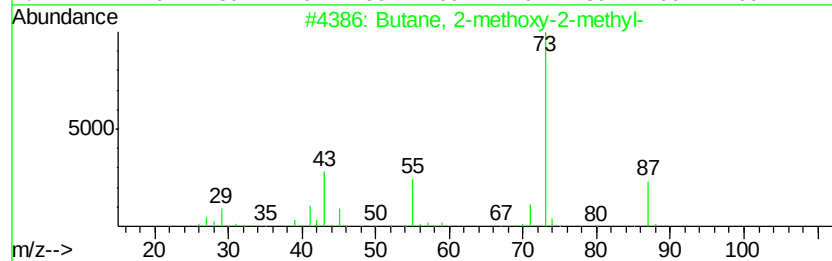
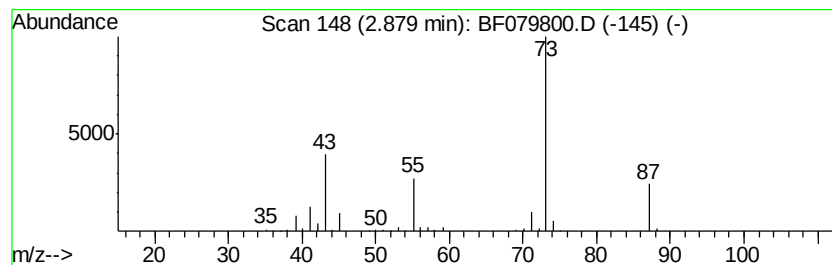
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TIC Integration Parameters: LSCINT.P

Peak Number 3 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.88	19.54 ng	335412	1,4-Dichlorobenzene-d4	7.47

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	10
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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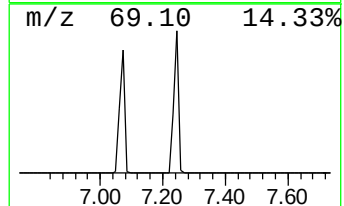
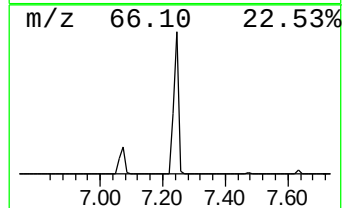
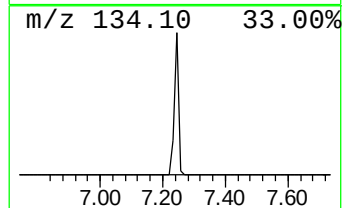
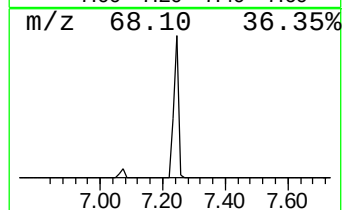
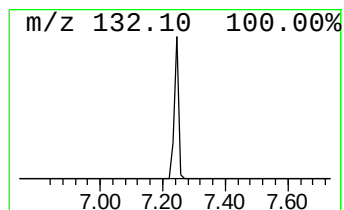
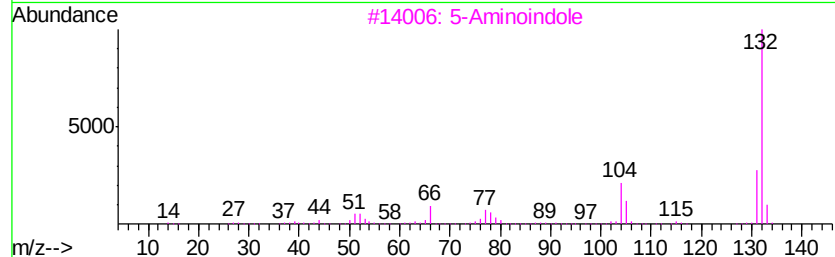
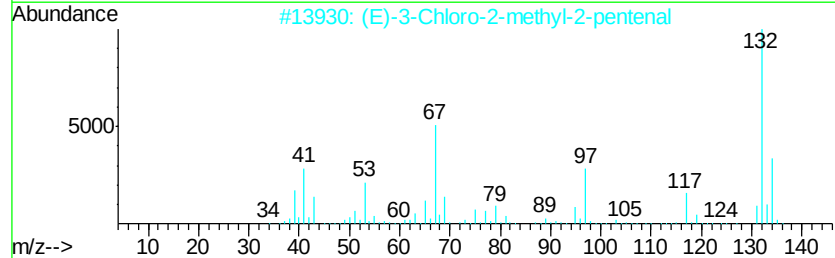
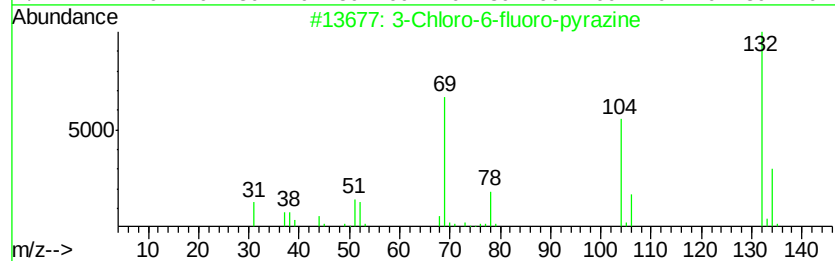
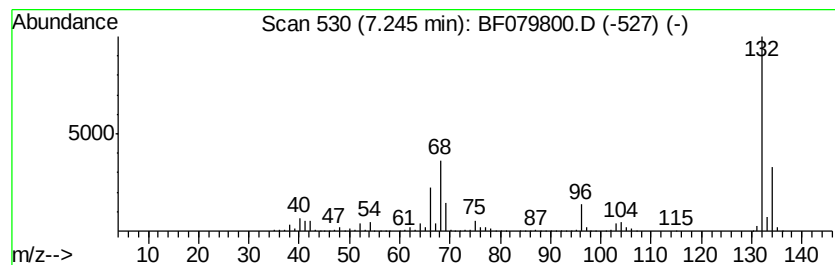
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Peak Number 4 unknown7.24 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.24	114.42 ng	1963820	1,4-Dichlorobenzene-d4	7.47

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	25
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 2,2-dime...	2.42	390.2	ng	6696630	1	7.47	343270	20.0
Butane, 2-methoxy...	2.88	19.5	ng	335412	1	7.47	343270	20.0
unknown7.24	7.24	114.4	ng	1963820	1	7.47	343270	20.0