

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF111017\
 Data File : BF100422.D
 Acq On : 11 Nov 2017 5:43
 Operator : SJ/JU
 Sample : I6355-12
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 110317-SIDI-E-D-S

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_F\METHODS\8270-BF110817.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	2.210	17	21	33	rVB2	88385	148904	3.34%	0.545%
2	5.122	512	516	525	rVB	281289	293169	6.58%	1.073%
3	5.516	579	583	586	rBV	1800816	1543285	34.62%	5.651%
4	6.516	749	753	756	rBV	1218556	1002690	22.49%	3.671%
5	6.545	756	758	764	rVB	64989	54939	1.23%	0.201%
6	6.669	774	779	782	rBV	2862978	2556341	57.35%	9.360%
7	6.898	815	818	822	rVB	1088131	977760	21.94%	3.580%
8	7.057	841	845	849	rVB	3223669	3265126	73.25%	11.955%
9	7.469	909	915	917	rBV	2492569	2798100	62.77%	10.245%
10	8.181	1032	1036	1040	rBV	1380923	1203150	26.99%	4.405%
11	9.263	1214	1220	1223	rBV	4309798	4457496	100.00%	16.321%
12	9.645	1282	1285	1292	rBV	79969	69928	1.57%	0.256%
13	9.939	1329	1335	1343	rVB	1350125	1229217	27.58%	4.501%
14	10.604	1445	1448	1452	rVB	135276	99210	2.23%	0.363%
15	10.645	1452	1455	1460	rVB	85912	80887	1.81%	0.296%
16	10.722	1464	1468	1472	rBV	1620515	1521006	34.12%	5.569%
17	11.422	1584	1587	1591	rBV	1195523	999364	22.42%	3.659%
18	13.010	1852	1857	1860	rBV	3487755	3435124	77.06%	12.577%
19	14.063	2032	2036	2039	rBV	950148	895372	20.09%	3.278%
20	15.545	2283	2288	2299	rBV	504867	680788	15.27%	2.493%

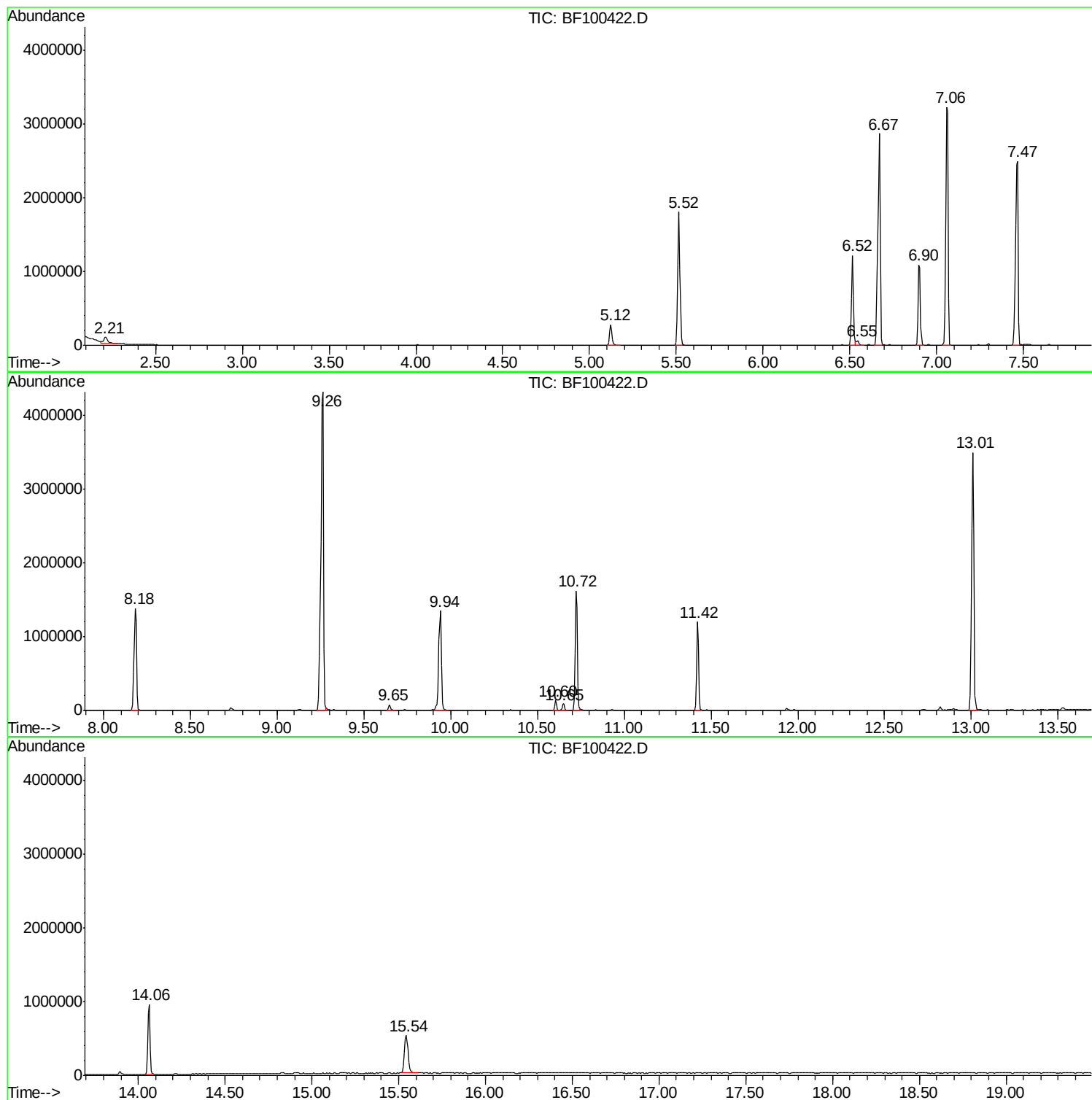
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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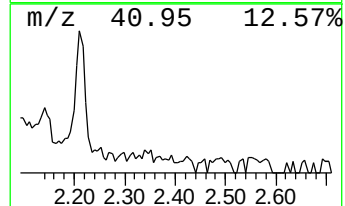
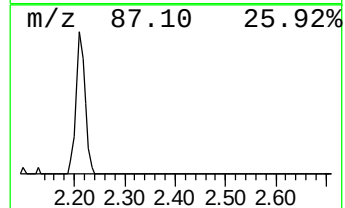
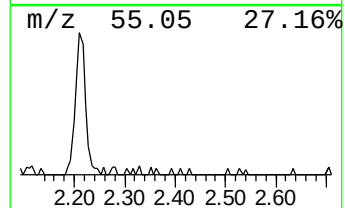
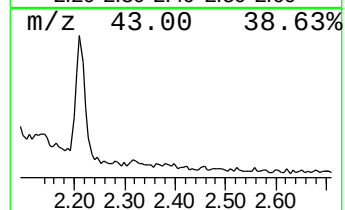
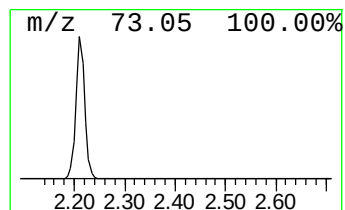
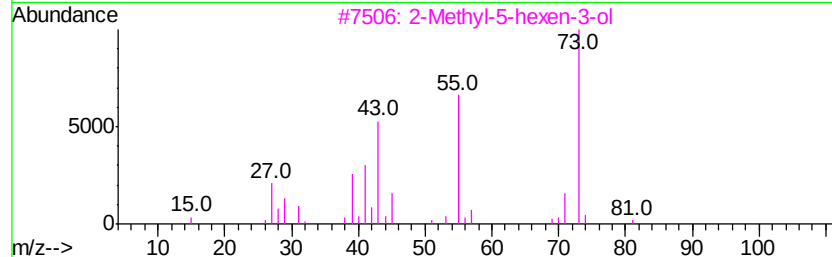
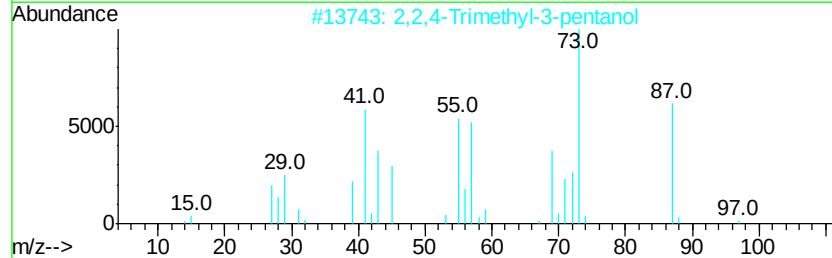
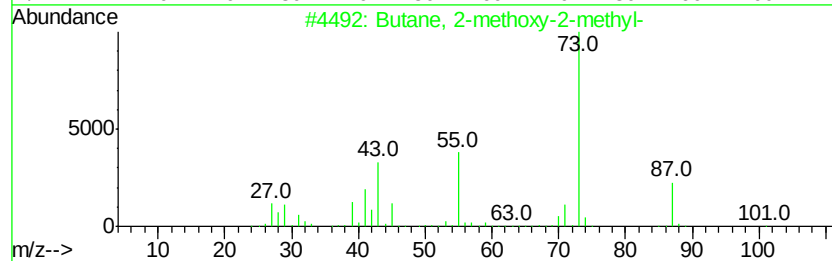
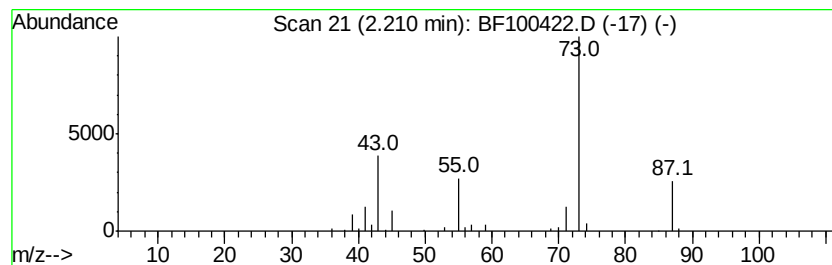
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
2.21	3.05 ng	148904	1,4-Dichlorobenzene-d4	6.90

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	39
3			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	25
4			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9



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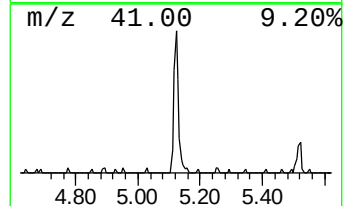
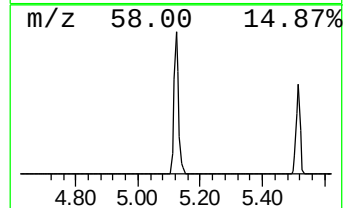
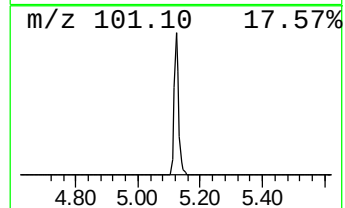
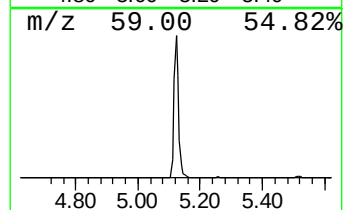
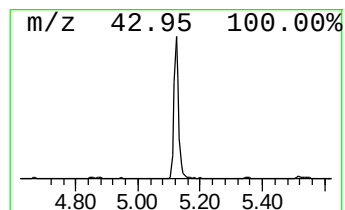
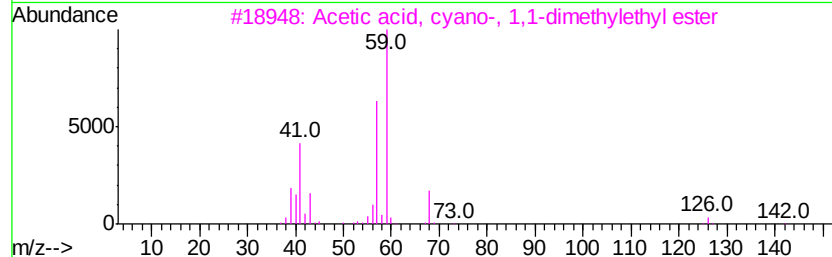
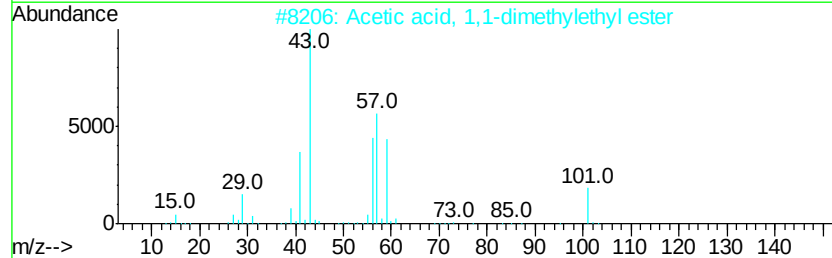
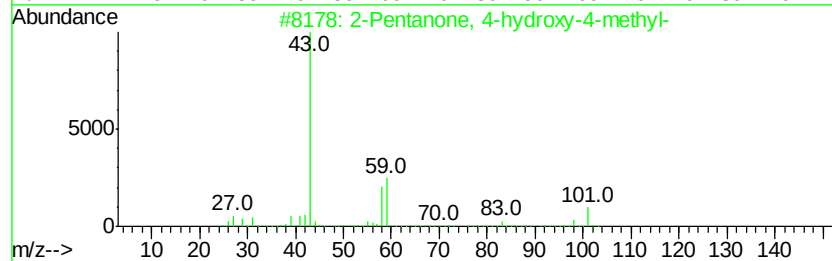
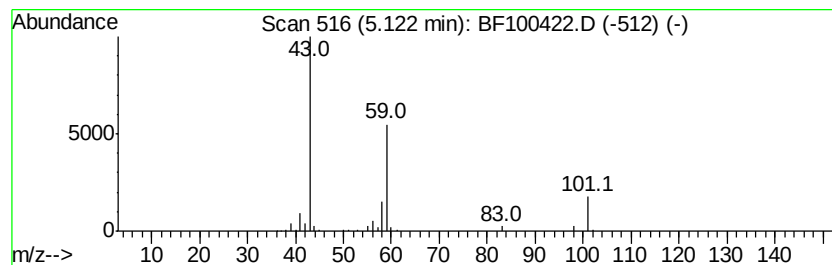
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	6.00 ng	293169	1,4-Dichlorobenzene-d4	6.90

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	32
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			3-Hexanol, 4-ethyl-	130	C8H18O	019780-44-0	9



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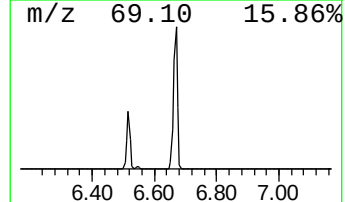
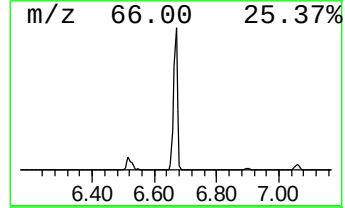
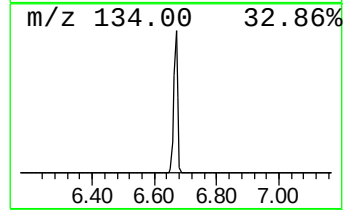
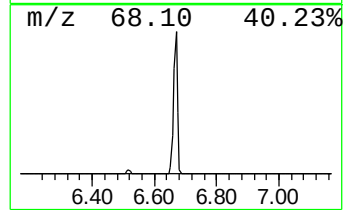
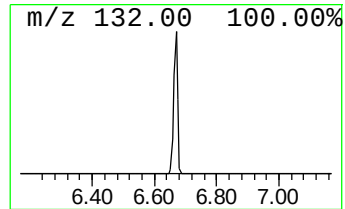
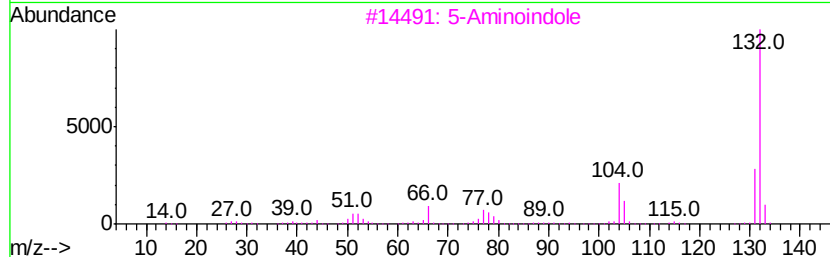
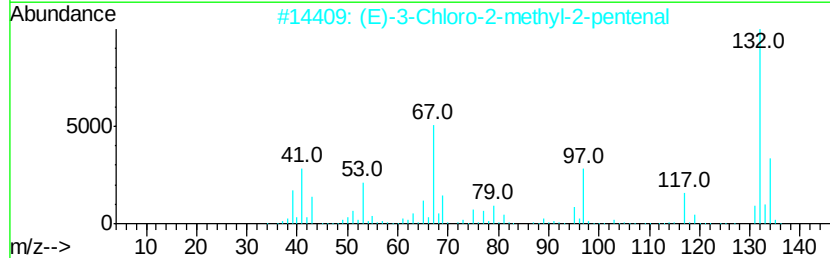
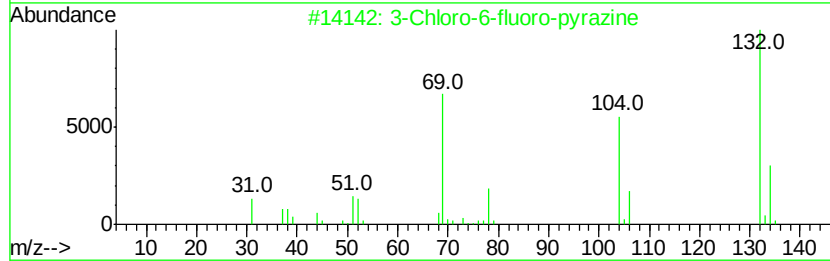
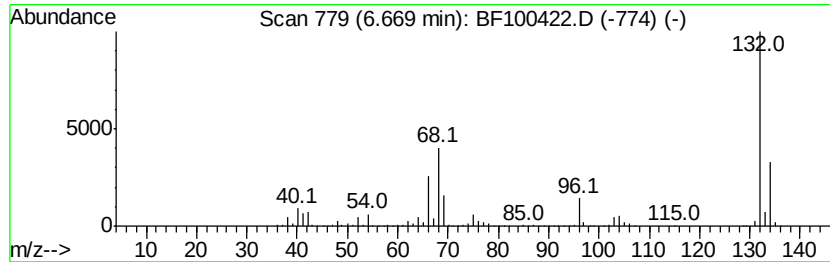
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Peak Number 3 unknown6.67 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.67	52.29 ng	2556340	1,4-Dichlorobenzene-d4	6.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.21	3.0	ng	148904	1	6.90	977760	20.0
2-Pentanone, 4-hy...	5.12	6.0	ng	293169	1	6.90	977760	20.0
unknown6.67	6.67	52.3	ng	2556340	1	6.90	977760	20.0