

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF111819\
 Data File : BF117832.D
 Acq On : 18 Nov 2019 13:48
 Operator : JU
 Sample : PB124875BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB124875BL

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.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	2.322	28	40	45	rVB	86318	123774	1.35%	0.179%
2	5.157	515	522	533	rBV	1086002	1454213	15.92%	2.108%
3	5.546	582	588	591	rBV	6357324	6644076	72.72%	9.631%
4	6.545	751	758	760	rBV	6023797	6870069	75.19%	9.959%
5	6.692	777	783	786	rBV	6474834	7119342	77.92%	10.320%
6	6.922	818	822	825	rBV	1781026	1582740	17.32%	2.294%
7	7.081	844	849	852	rBV	5779384	5727051	62.68%	8.302%
8	7.481	911	917	920	rBV	4209483	4531447	49.60%	6.569%
9	8.204	1035	1040	1044	rBV	2177498	2033267	22.25%	2.947%
10	9.286	1218	1224	1227	rBV	8457213	8391102	91.84%	12.164%
11	9.969	1334	1340	1343	rBV	2890658	2407448	26.35%	3.490%
12	10.763	1468	1475	1478	rBV	5743172	5837942	63.90%	8.463%
13	11.463	1587	1594	1597	rBV	2890215	2620067	28.68%	3.798%
14	13.057	1860	1865	1868	rBV	9173573	9136586	100.00%	13.244%
15	14.110	2040	2044	2054	rVB	2557942	2354877	25.77%	3.414%
16	15.615	2294	2300	2316	rBV	1606468	2150693	23.54%	3.118%

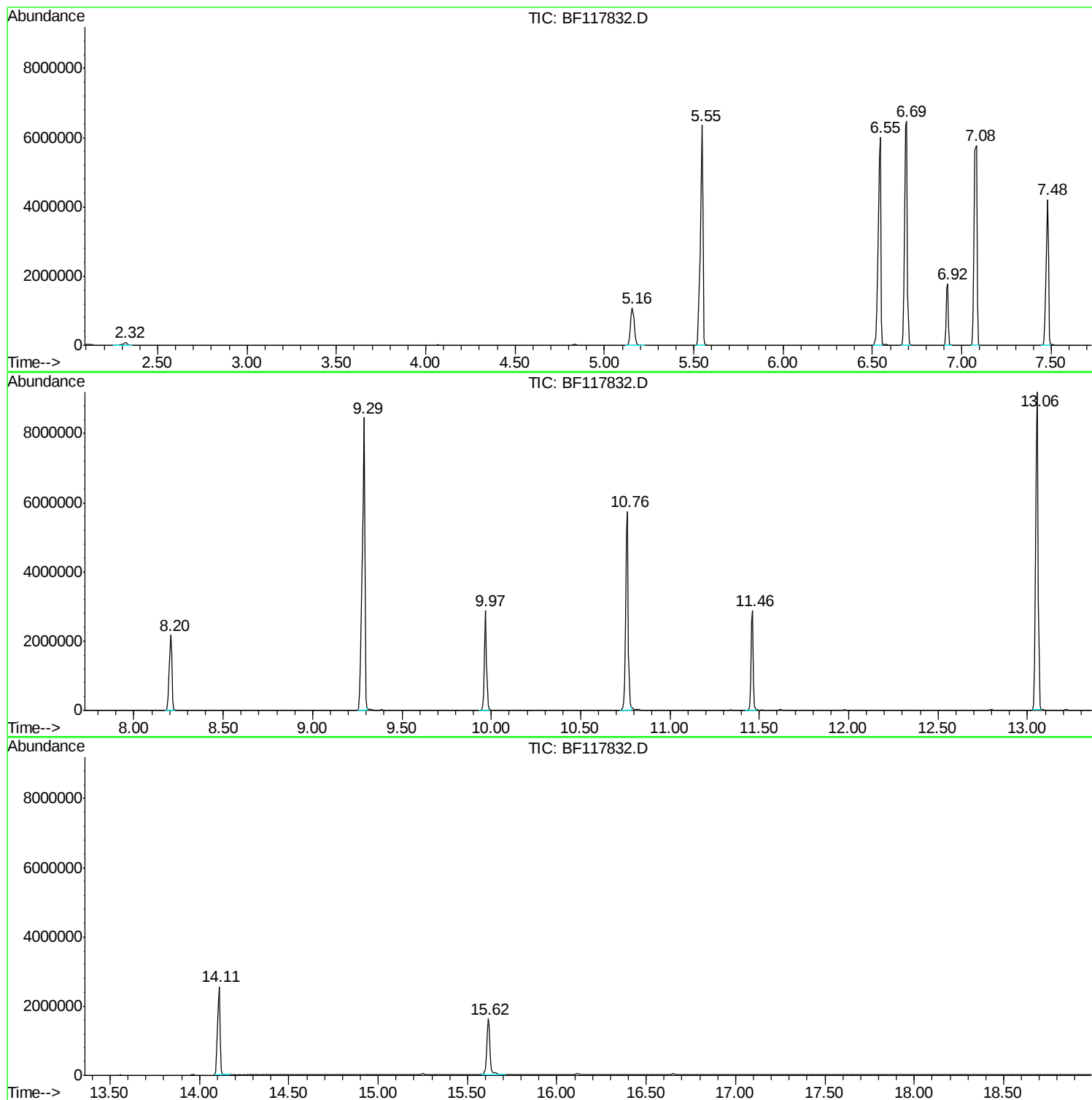
Sum of corrected areas: 68984694

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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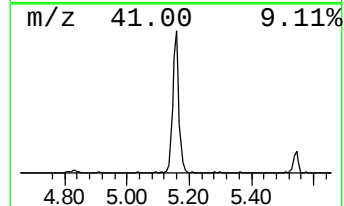
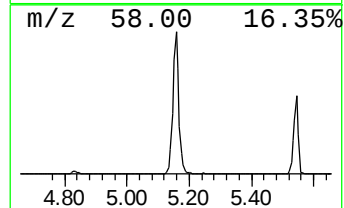
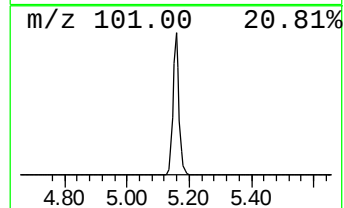
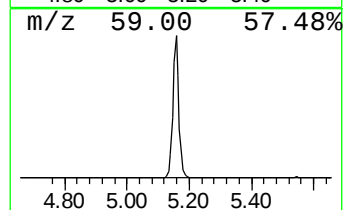
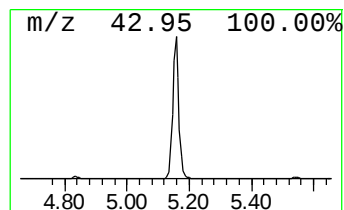
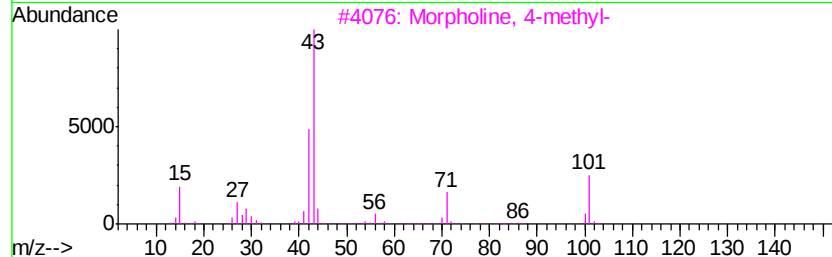
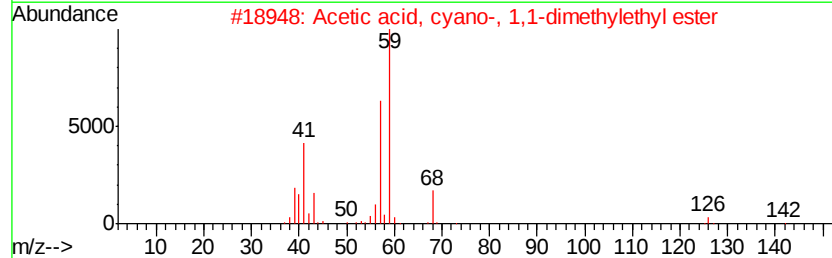
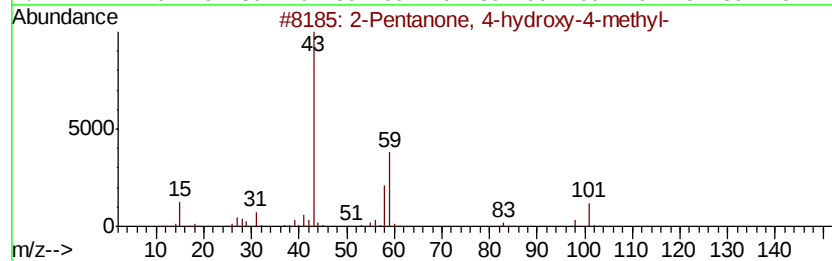
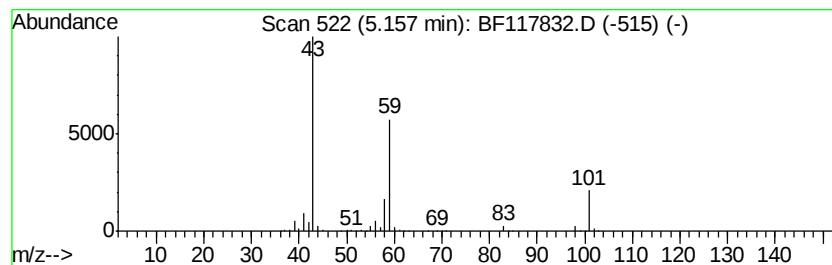
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	18.38 ng	1454210	1,4-Dichlorobenzene-d4	6.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4		2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9
5		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	9



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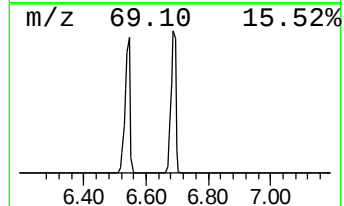
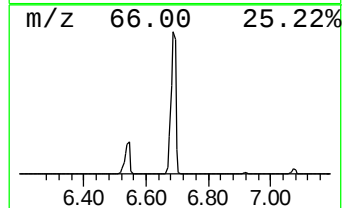
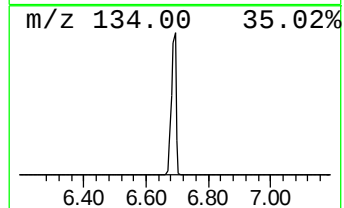
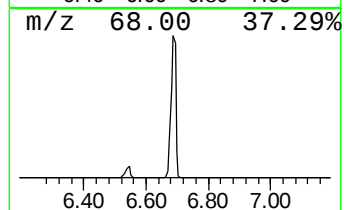
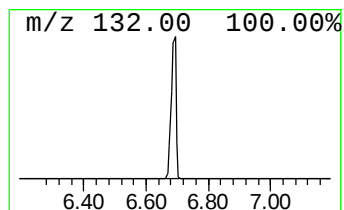
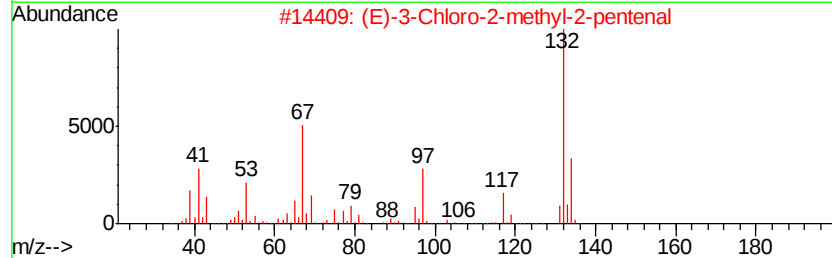
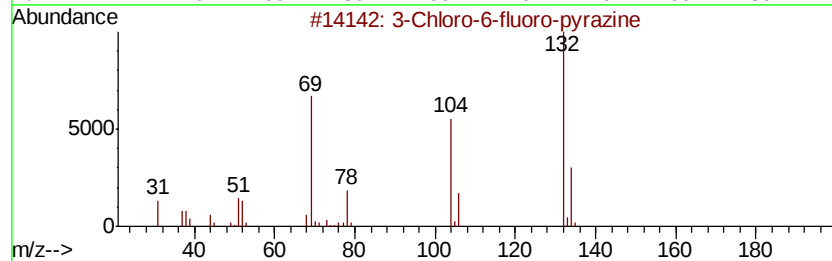
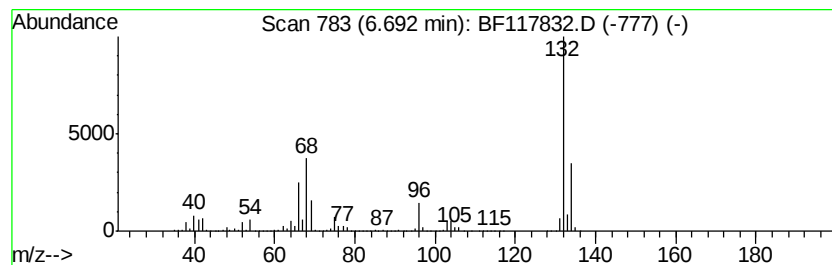
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Peak Number 2 unknown6.69 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.69	89.96 ng	7119340	1,4-Dichlorobenzene-d4	6.92

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		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	16
4		5-Aminoindole	132	C8H8N2	005192-03-0	16
5		Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.16	18.4	ng	1454210	1	6.92	1582740	20.0
unknown6.69	6.69	90.0	ng	7119340	1	6.92	1582740	20.0