

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF070318\
 Data File : BF107087.D
 Acq On : 3 Jul 2018 17:34
 Operator : JU/SJ
 Sample : J3245-19 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-01-062918-B

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.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	5.178	479	483	489	rBV2	62222	69786	9.47%	1.061%
2	5.596	551	554	570	rBV	516134	542449	73.58%	8.244%
3	6.484	701	705	709	rBV	23072	20547	2.79%	0.312%
4	6.601	722	725	735	rBV	500538	546788	74.17%	8.310%
5	6.737	744	748	752	rBV	583646	542930	73.64%	8.251%
6	6.954	782	785	789	rBV	381499	309010	41.91%	4.696%
7	7.113	808	812	817	rVB	508138	427474	57.98%	6.497%
8	7.519	877	881	891	rBV	321960	288006	39.06%	4.377%
9	8.242	1000	1004	1014	rBV	343226	359951	48.82%	5.470%
10	9.313	1182	1186	1194	rBV	585142	579300	78.58%	8.804%
11	9.707	1248	1253	1259	rBV	56761	71010	9.63%	1.079%
12	9.854	1275	1278	1283	rVV2	6072	9561	1.30%	0.145%
13	9.901	1283	1286	1290	rVB2	9705	9171	1.24%	0.139%
14	10.001	1299	1303	1315	rVB	393761	380706	51.64%	5.786%
15	10.795	1435	1438	1446	rBV	343372	331056	44.90%	5.031%
16	11.495	1553	1557	1560	rBV	477806	421203	57.13%	6.401%
17	12.113	1658	1662	1664	rBV5	9872	11814	1.60%	0.180%
18	12.495	1725	1727	1729	rBV2	12257	9532	1.29%	0.145%
19	12.548	1733	1736	1739	rVV3	12425	12924	1.75%	0.196%
20	12.713	1761	1764	1768	rBV	55188	52579	7.13%	0.799%
21	12.895	1792	1795	1798	rBV4	12519	11937	1.62%	0.181%
22	12.924	1798	1800	1801	rVV2	12817	12137	1.65%	0.184%
23	12.948	1801	1804	1808	rVB	64062	62395	8.46%	0.948%
24	13.077	1822	1826	1829	rBV	835901	737250	100.00%	11.204%
25	13.277	1854	1860	1862	rBV6	15520	29858	4.05%	0.454%
26	13.377	1875	1877	1880	rBV2	10025	11938	1.62%	0.181%
27	13.471	1891	1893	1896	rBV	14820	14747	2.00%	0.224%
28	13.954	1972	1975	1978	rBV2	38404	37639	5.11%	0.572%
29	14.148	2004	2008	2011	rBV	413617	408588	55.42%	6.209%
30	15.695	2267	2271	2279	rVB	192979	257777	34.96%	3.918%

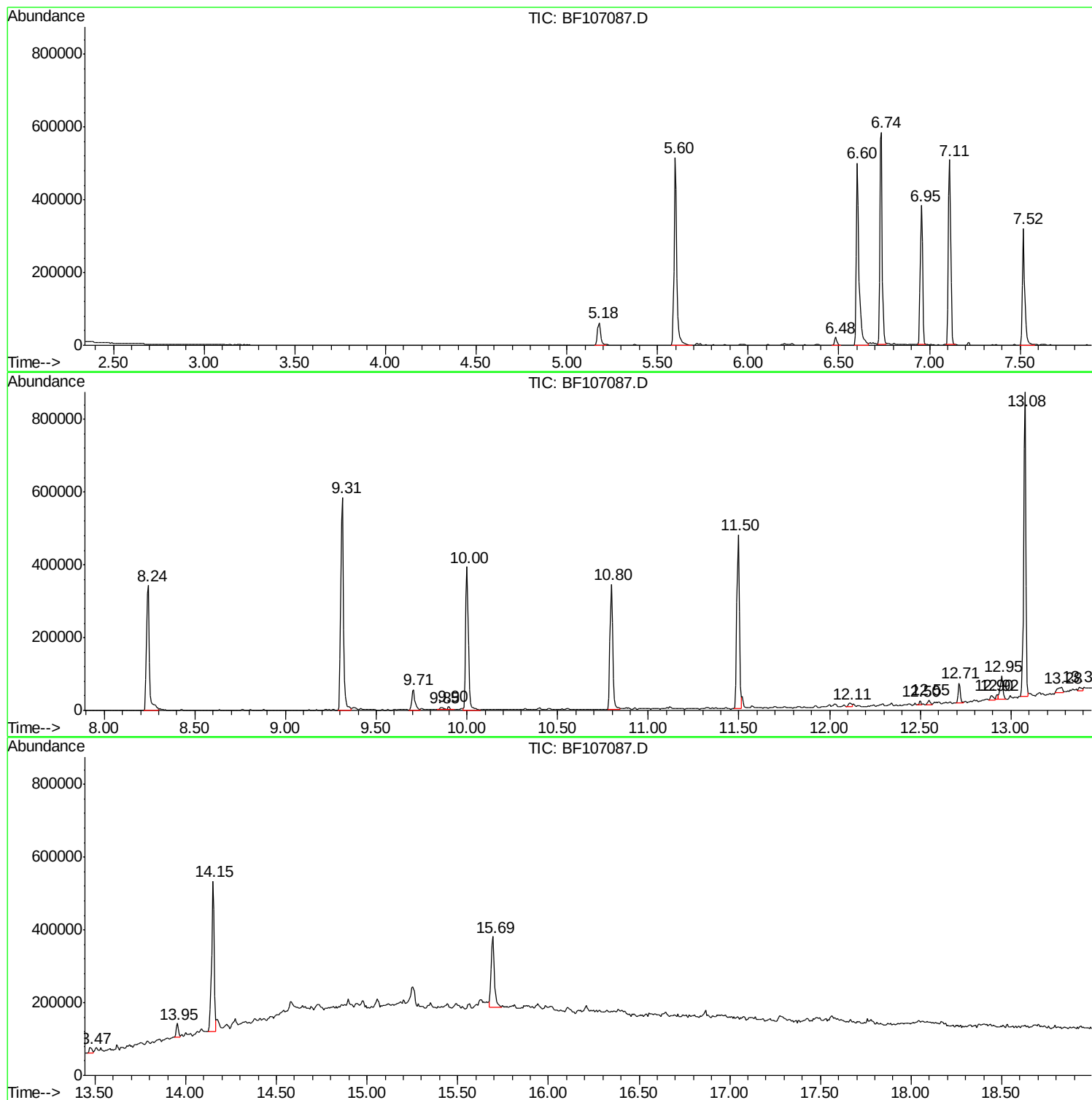
Sum of corrected areas: 6580063

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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.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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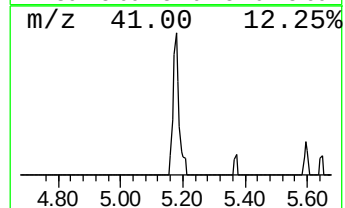
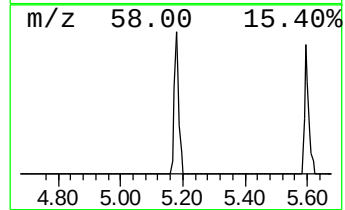
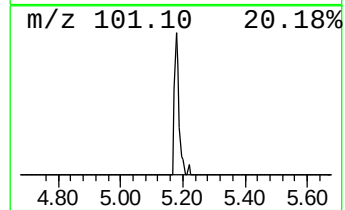
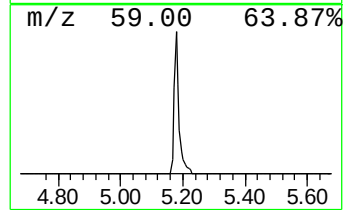
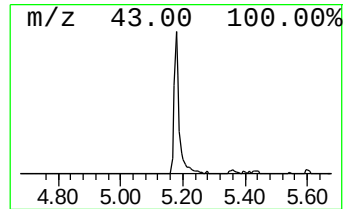
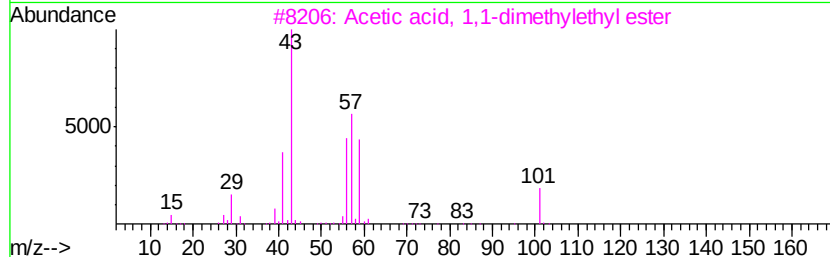
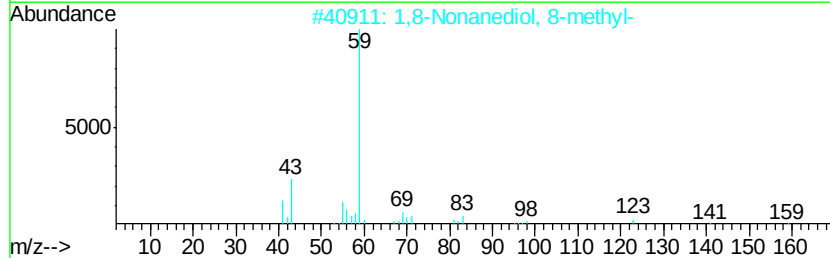
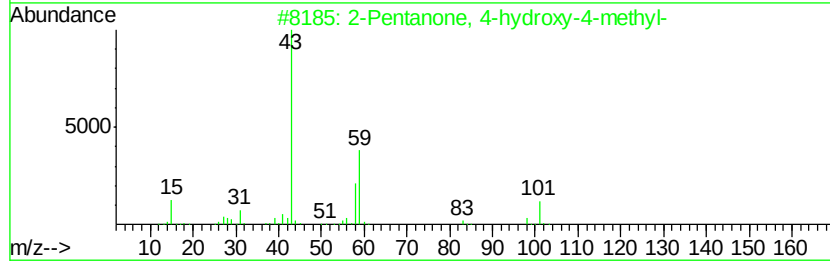
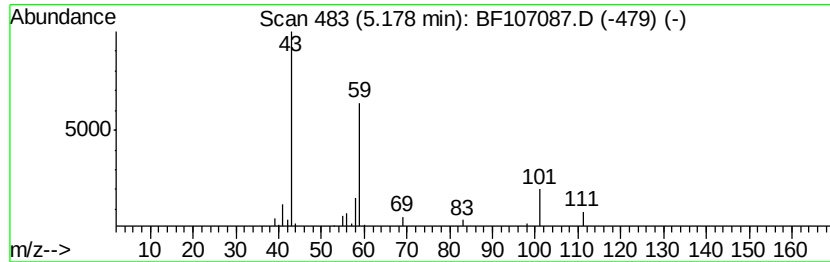
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.18	4.52 ng	69786	1,4-Dichlorobenzene-d4	6.95	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2	1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	37
3	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	36
4	n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	28
5	5-Hexen-2-one	98	C6H10O	000109-49-9	9



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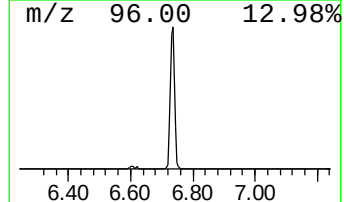
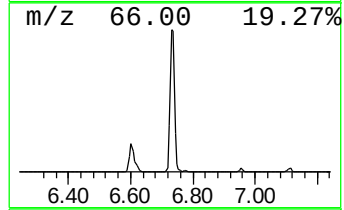
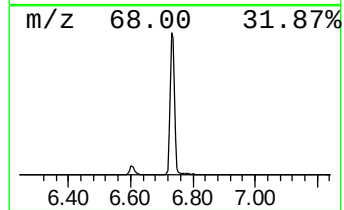
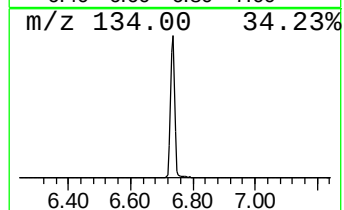
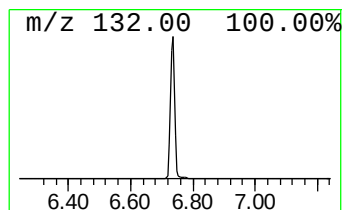
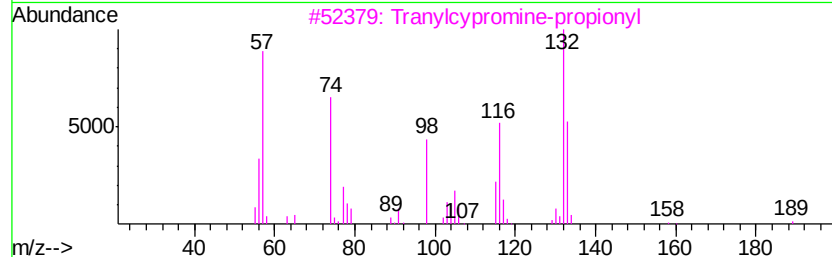
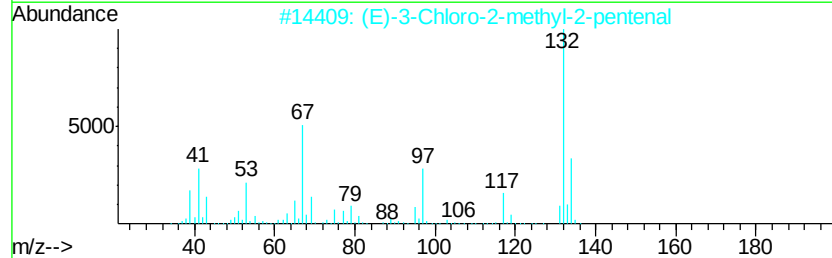
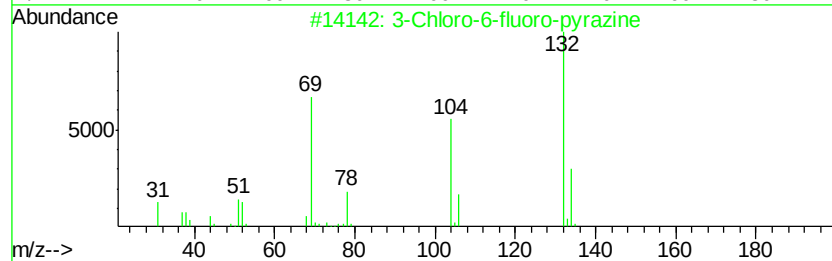
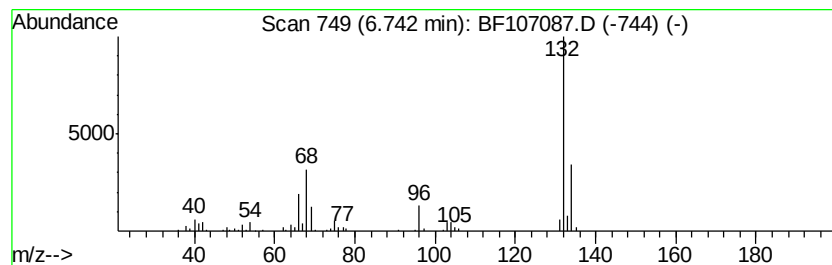
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Peak Number 2 unknown6.74 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	35.14 ng	542930	1,4-Dichlorobenzene-d4	6.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	12
4		Xenon	132	Xe	007440-63-3	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.18	4.5	ng	69786	1	6.95	309010	20.0
unknown6.74	6.74	35.1	ng	542930	1	6.95	309010	20.0