

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061721\
 Data File : BF124361.D
 Acq On : 17 Jun 2021 14:07
 Operator : JU/CG
 Sample : M2746-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NEW-DORP-2

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-BF060321.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF124361.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.998	492	495	501	rBV	26200	31485	2.57%	0.356%
2	5.410	561	565	581	rBV	1102513	986869	80.47%	11.161%
3	6.434	734	739	749	rBV	1017317	976707	79.64%	11.046%
4	6.781	795	798	802	rVB	285642	224780	18.33%	2.542%
5	7.351	891	895	898	rBV	807216	660524	53.86%	7.470%
6	7.975	998	1001	1008	rBV	49154	52250	4.26%	0.591%
7	8.069	1013	1017	1020	rBV	342400	284082	23.16%	3.213%
8	9.145	1196	1200	1203	rBV	1663312	1226413	100.00%	13.870%
9	9.822	1312	1315	1318	rBV	461852	334222	27.25%	3.780%
10	10.616	1446	1450	1453	rBV	981053	805663	65.69%	9.112%
11	11.310	1565	1568	1571	rBV	422935	337878	27.55%	3.821%
12	11.333	1571	1572	1575	rVB	54085	30088	2.45%	0.340%
13	11.839	1655	1658	1661	rBV2	29924	27526	2.24%	0.311%
14	11.921	1670	1672	1676	rBV2	18564	20551	1.68%	0.232%
15	12.527	1771	1775	1778	rBV	203966	178307	14.54%	2.017%
16	12.757	1807	1814	1817	rBV	169970	182004	14.84%	2.058%
17	12.898	1832	1838	1841	rBV	1250631	1046098	85.30%	11.831%
18	12.974	1847	1851	1855	rBV4	10285	16949	1.38%	0.192%
19	13.057	1860	1865	1868	rBV4	11069	18138	1.48%	0.205%
20	13.086	1868	1870	1872	rVB2	30058	22761	1.86%	0.257%
21	13.151	1878	1881	1884	rVB2	24579	21126	1.72%	0.239%
22	13.186	1884	1887	1889	rBV2	15192	18080	1.47%	0.204%
23	13.598	1954	1957	1962	rVB	21092	24539	2.00%	0.278%
24	13.704	1972	1975	1978	rBV	21077	18329	1.49%	0.207%
25	13.763	1982	1985	1988	rBV2	16340	21434	1.75%	0.242%
26	13.810	1989	1993	1996	rBV3	21030	25654	2.09%	0.290%
27	13.957	2011	2018	2021	rBV2	364663	419666	34.22%	4.746%
28	13.980	2021	2022	2030	rVB	112021	77880	6.35%	0.881%
29	14.062	2033	2036	2038	rVB2	21326	17728	1.45%	0.200%
30	14.110	2041	2044	2048	rBV4	17253	25264	2.06%	0.286%
31	14.345	2079	2084	2086	rBV	20918	18637	1.52%	0.211%
32	14.992	2190	2194	2197	rBV	108334	140792	11.48%	1.592%
33	15.098	2209	2212	2218	rBV	26630	35790	2.92%	0.405%
34	15.292	2241	2245	2251	rVB	52404	66923	5.46%	0.757%
35	15.357	2251	2256	2261	rBV	83799	104488	8.52%	1.182%
36	15.415	2261	2266	2270	rBV	175939	233841	19.07%	2.645%

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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

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37	15.851	2335	2340	2344	rBV5	10364	15862	1.29%	0.179%
38	16.862	2507	2512	2517	rBV3	30644	54707	4.46%	0.619%
39	17.298	2582	2586	2592	rBV3	16957	38000	3.10%	0.430%

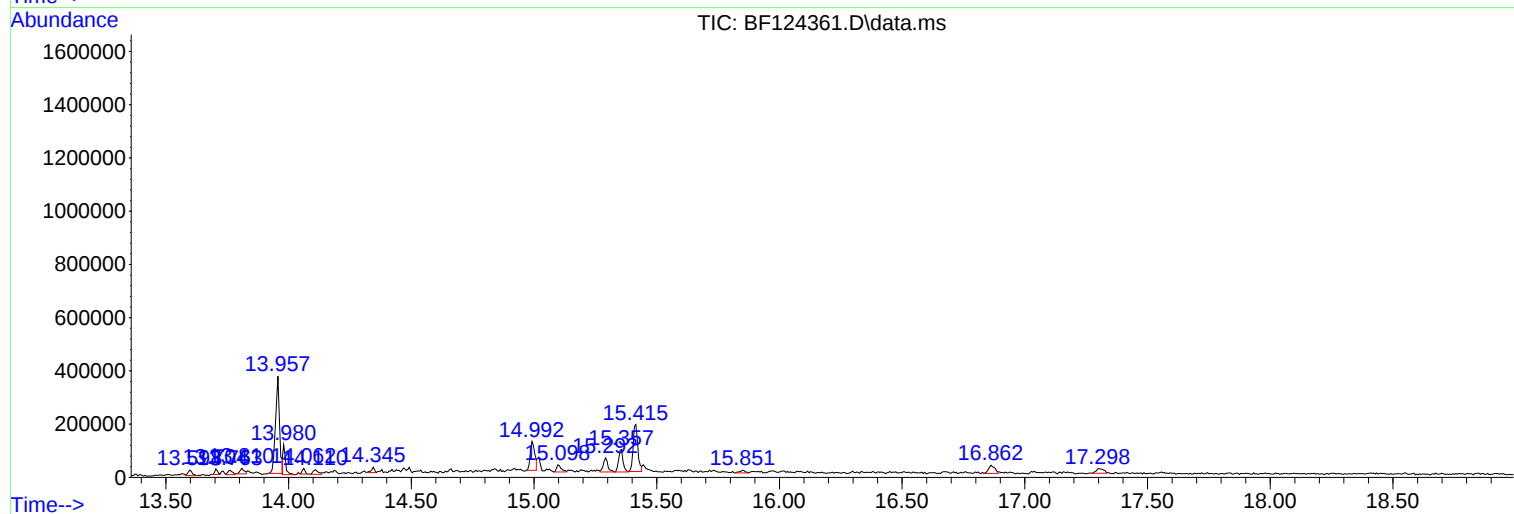
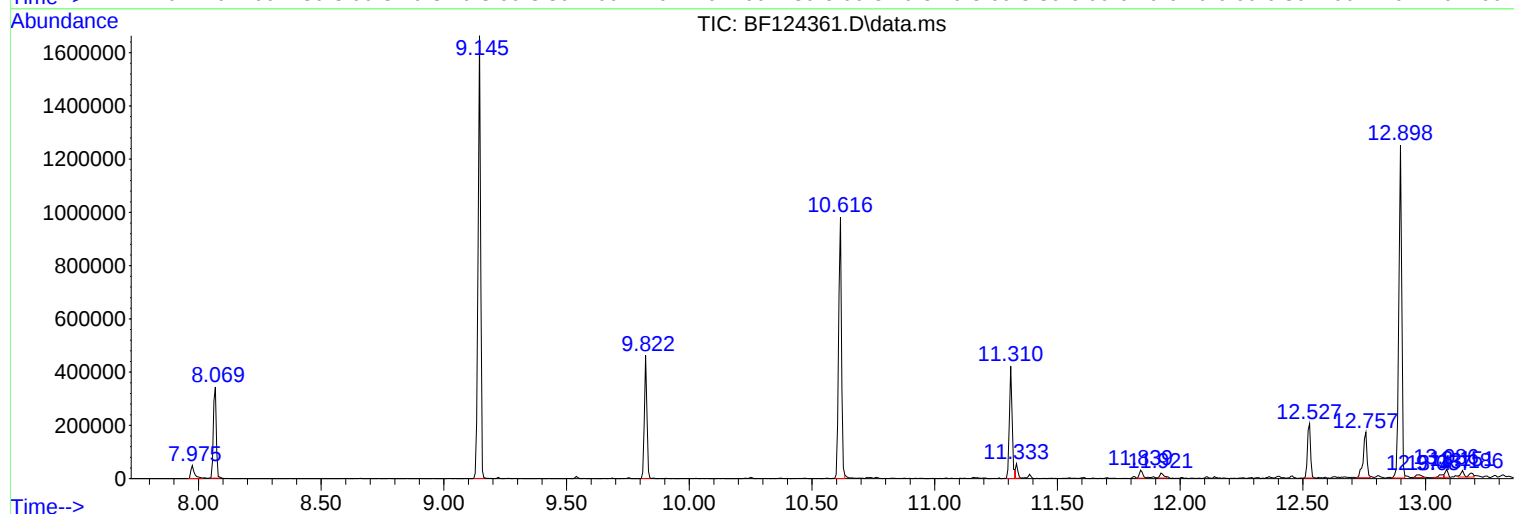
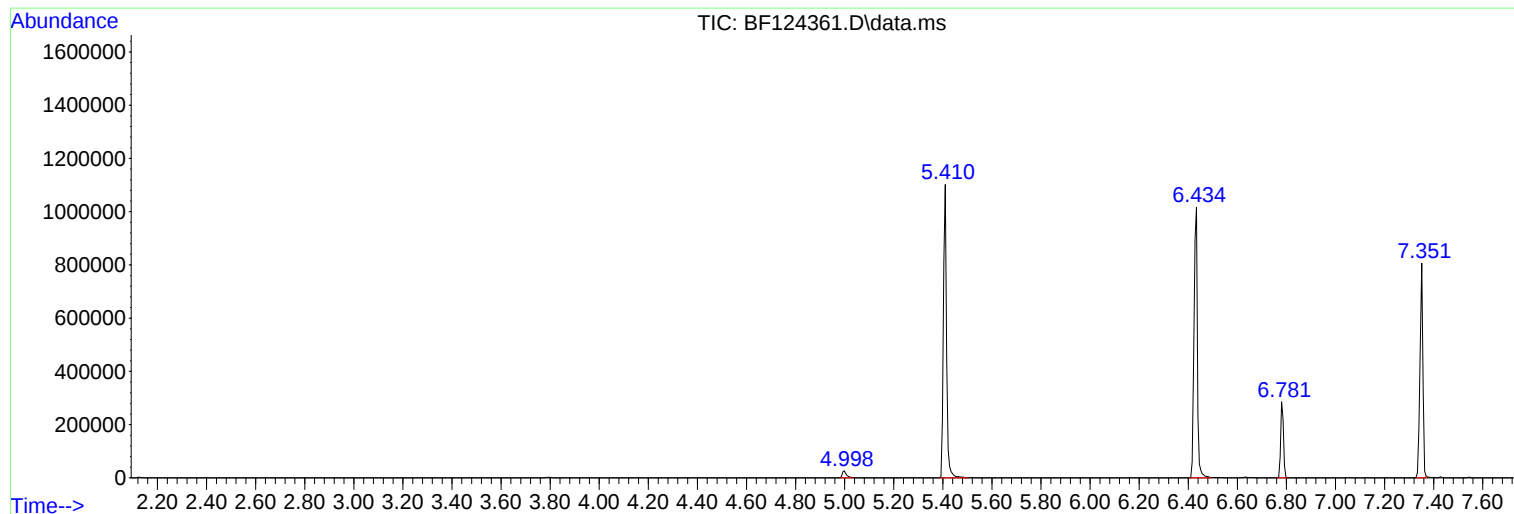
Sum of corrected areas: 8842035

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.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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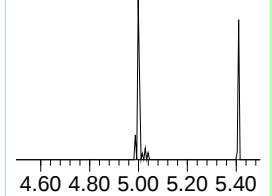
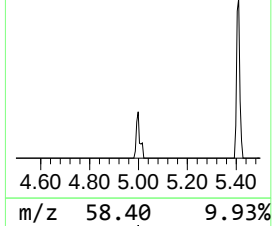
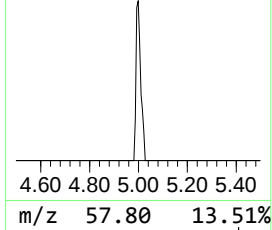
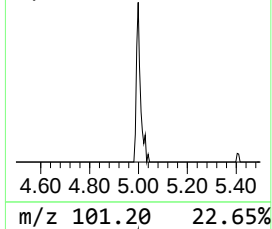
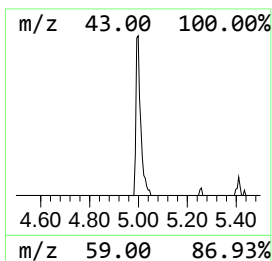
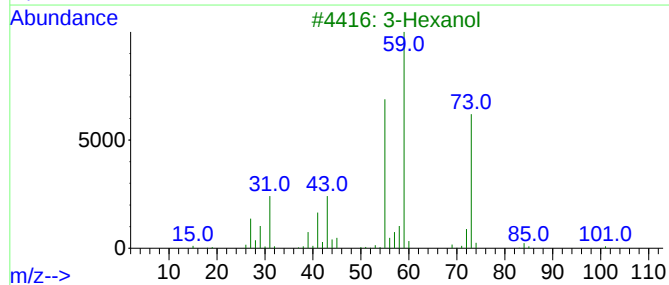
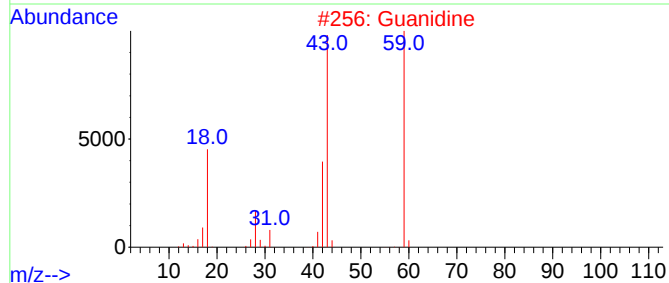
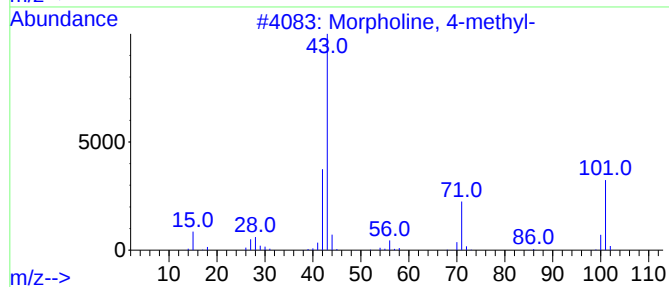
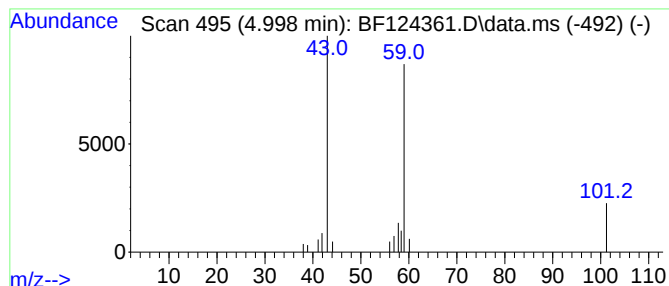
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown4.998 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.998	2.80 ng	31485	1,4-Dichlorobenzene-d4	6.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
2			Guanidine	59	CH5N3	000113-00-8	9
3			3-Hexanol	102	C6H14O	000623-37-0	9
4			2-Nonanone, 9-hydroxy-	158	C9H18O2	025368-56-3	9
5			Silane, ethyldimethyl-	88	C4H12Si	000758-21-4	9



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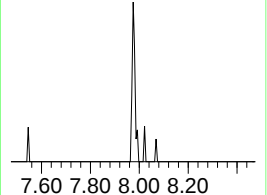
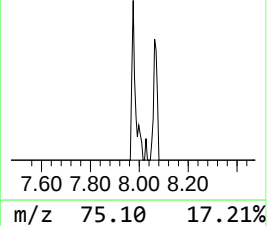
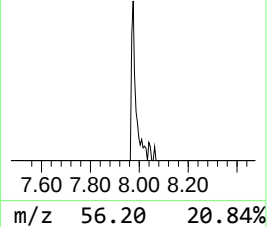
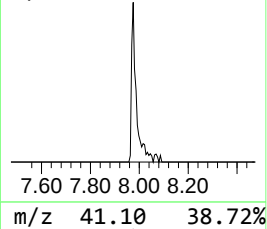
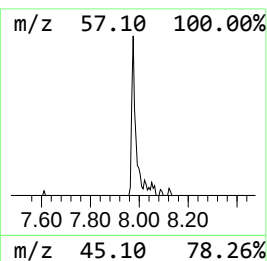
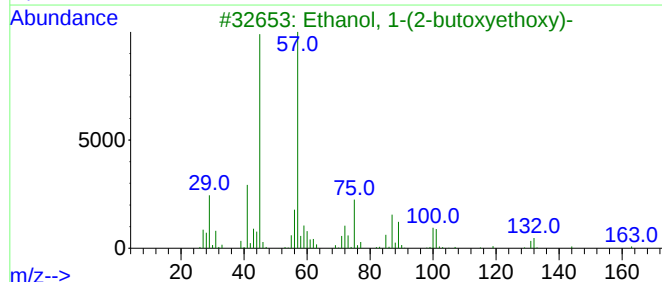
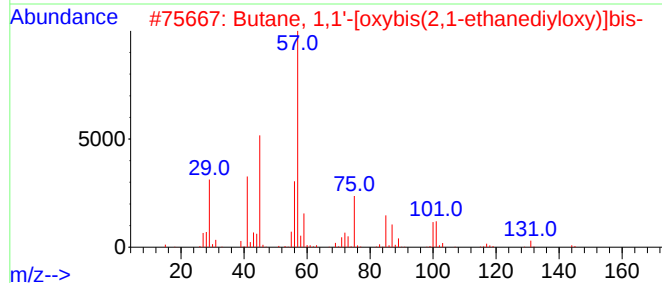
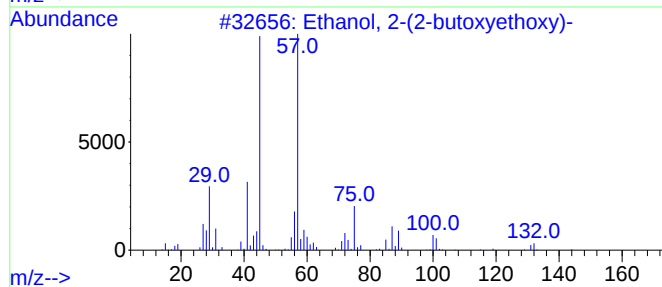
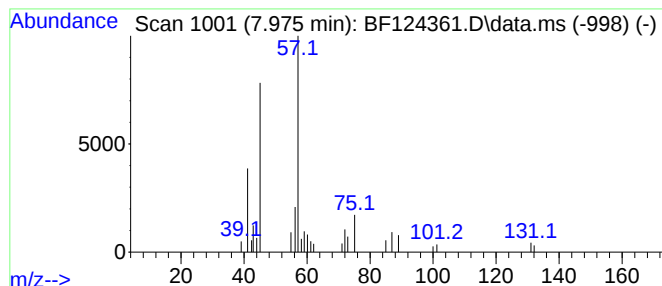
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Peak Number 2 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.975	3.68 ng	52250	Naphthalene-d8	8.069

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72
2			Butane, 1,1'-[oxybis(2,1-ethaned...	218	C12H26O3	000112-73-2	53
3			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	45
4			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	42
5			Ethanol, 2-[2-(2-butoxyethoxy)et...	206	C10H22O4	000143-22-6	42



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown4.998	4.998	2.8	ng	31485	1	6.781	224780	20.0
Ethanol, 2-(2-b...	7.975	3.7	ng	52250	2	8.069	284082	20.0