

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080824\
 Data File : BF138874.D
 Acq On : 08 Aug 2024 19:33
 Operator : RC/JU
 Sample : P3484-01
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B

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-BF073024.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF138874.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.416	219	225	237	rBV2	3148	12994	1.98%	0.243%
2	5.051	498	503	520	rVB	49076	70897	10.81%	1.324%
3	5.469	569	574	596	rBB	450383	612859	93.48%	11.442%
4	6.481	741	746	765	rBV	486972	635001	96.86%	11.855%
5	6.681	776	780	792	rBV	6034	12729	1.94%	0.238%
6	6.839	802	807	812	rBV	247169	297978	45.45%	5.563%
7	7.404	898	903	917	rBV	320406	405201	61.81%	7.565%
8	7.657	943	946	952	rBB	6045	7734	1.18%	0.144%
9	8.122	1019	1025	1030	rBV	292448	370399	56.50%	6.915%
10	8.433	1074	1078	1098	rBV	24146	42850	6.54%	0.800%
11	9.192	1202	1207	1212	rBV	541358	655593	100.00%	12.239%
12	9.869	1317	1322	1333	rBV	290131	358542	54.69%	6.694%
13	9.963	1333	1338	1342	rVB	15953	22729	3.47%	0.424%
14	10.663	1452	1457	1471	rBV	202669	276998	42.25%	5.171%
15	11.357	1570	1575	1585	rBV	226577	288497	44.01%	5.386%
16	11.886	1656	1665	1670	rBV5	5553	10449	1.59%	0.195%
17	11.968	1676	1679	1687	rVB2	4232	7331	1.12%	0.137%
18	12.498	1765	1769	1775	rVB	6986	9088	1.39%	0.170%
19	12.568	1775	1781	1786	rBV	44800	55363	8.44%	1.034%
20	12.798	1813	1820	1826	rBV	57330	78628	11.99%	1.468%
21	12.933	1838	1843	1853	rVB	336072	432243	65.93%	8.070%
22	13.015	1853	1857	1862	rBV4	5898	10396	1.59%	0.194%
23	13.115	1868	1874	1880	rBV5	7551	18031	2.75%	0.337%
24	13.227	1889	1893	1901	rVB2	7992	12923	1.97%	0.241%
25	13.321	1905	1909	1912	rBV	5428	7273	1.11%	0.136%
26	13.639	1955	1963	1966	rBV5	4671	9568	1.46%	0.179%
27	13.751	1978	1982	1984	rBV3	4842	7718	1.18%	0.144%
28	13.857	1996	2000	2014	rVB3	8828	23638	3.61%	0.441%
29	13.998	2018	2024	2036	rVV2	192933	298261	45.49%	5.568%
30	14.451	2097	2101	2104	rBV4	7937	13012	1.98%	0.243%
31	15.033	2196	2200	2214	rVB	19836	55536	8.47%	1.037%
32	15.333	2248	2251	2257	rVB	9446	14738	2.25%	0.275%
33	15.398	2258	2262	2266	rVV	12575	21340	3.26%	0.398%
34	15.456	2267	2272	2282	rVB	123245	199900	30.49%	3.732%

Sum of corrected areas: 5356437

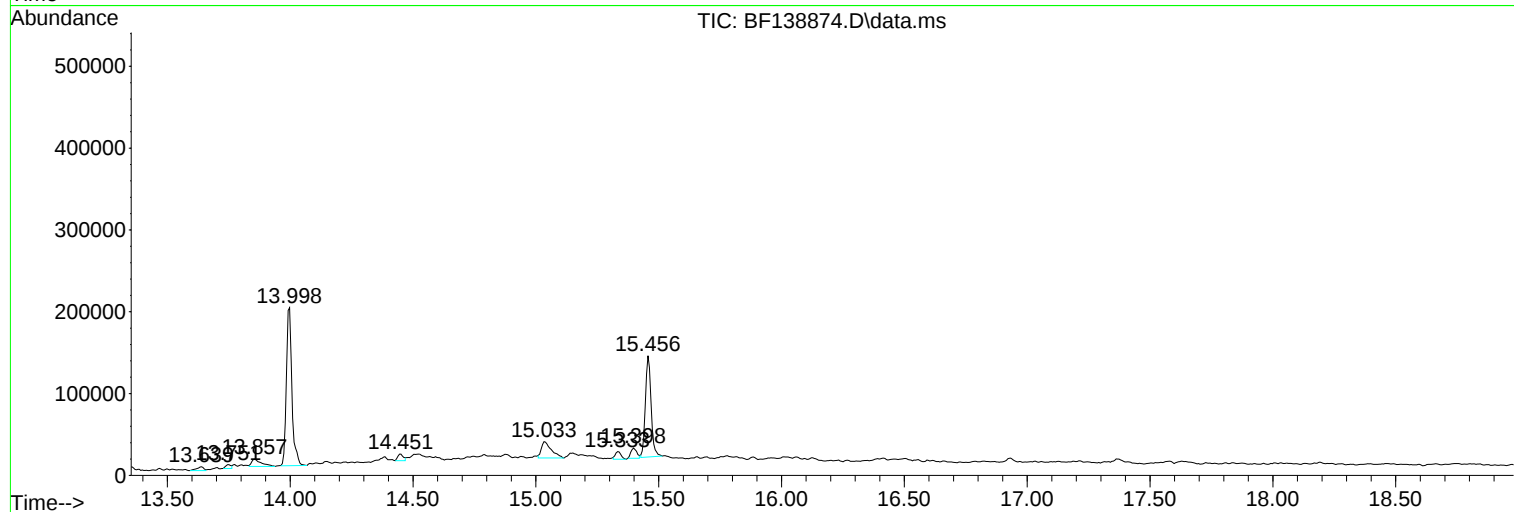
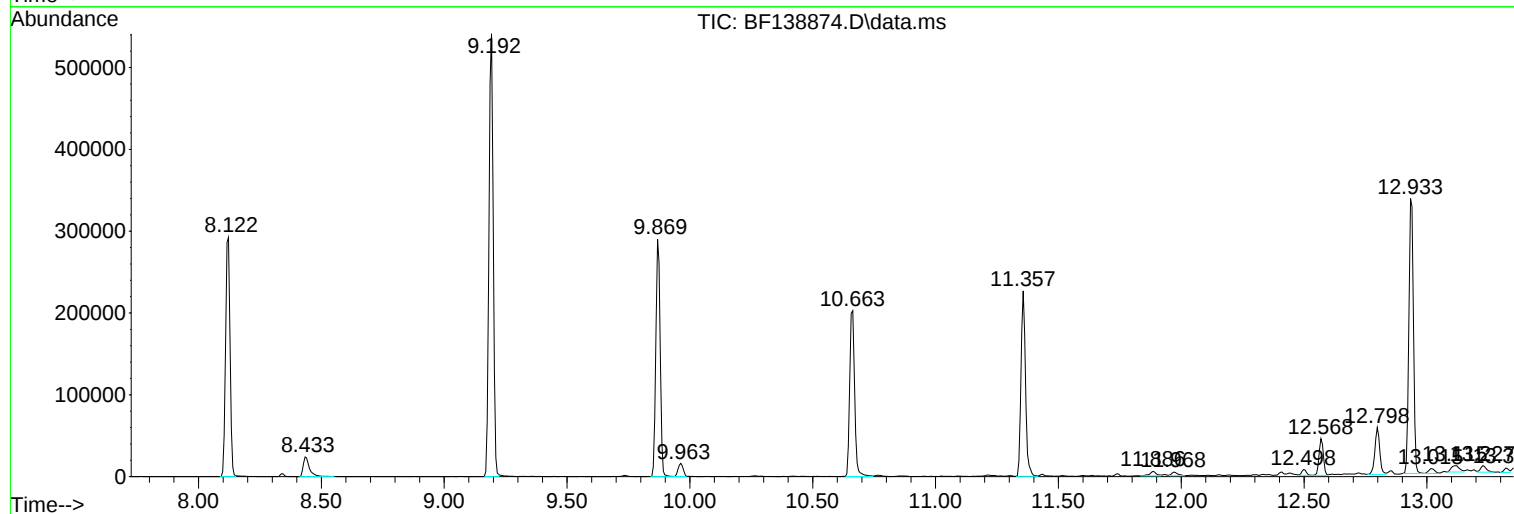
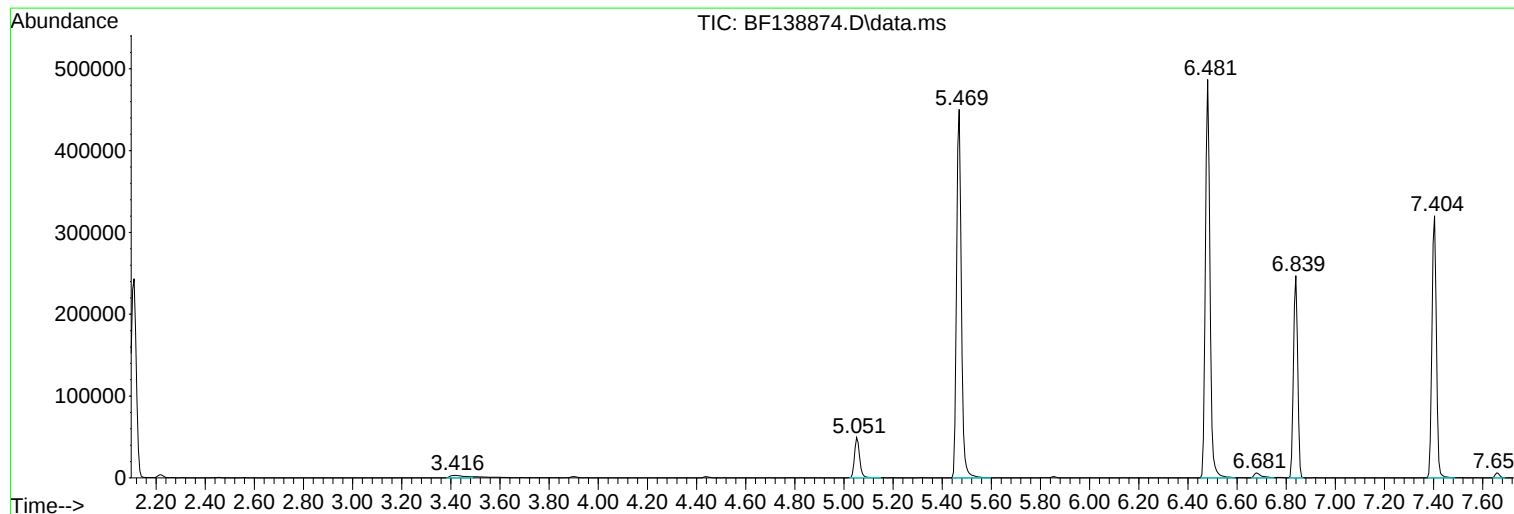
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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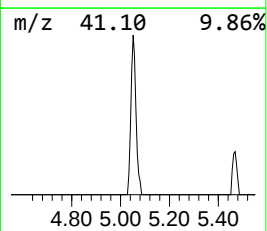
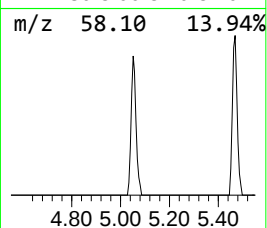
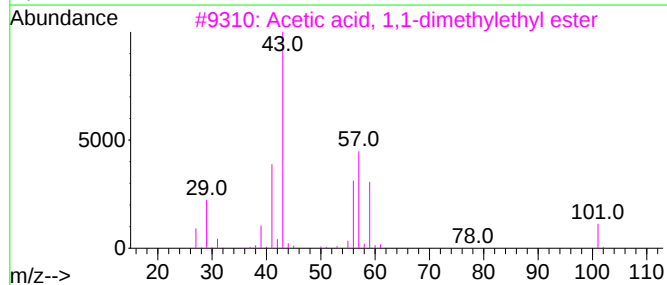
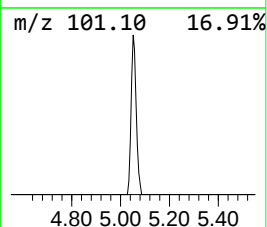
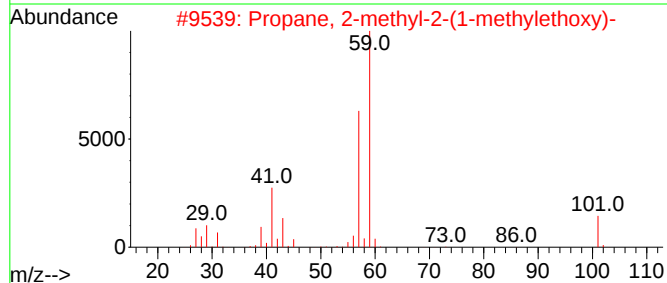
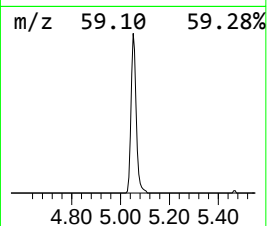
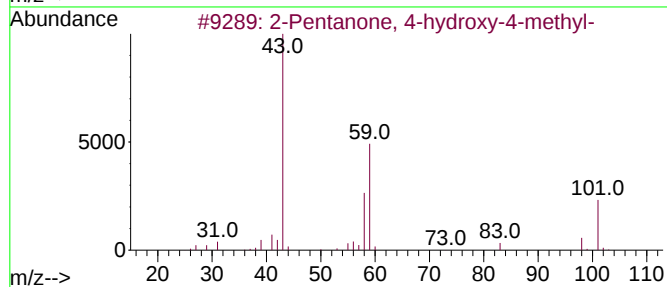
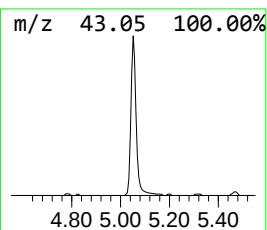
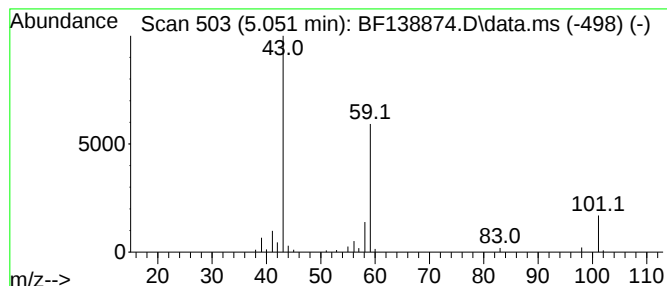
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.051	4.76 ng	70897	1,4-Dichlorobenzene-d4	6.839

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64		
2	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36		
3	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28		
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27		
5	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12		



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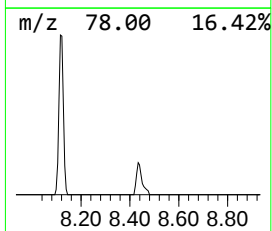
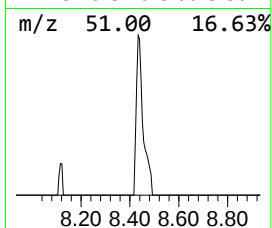
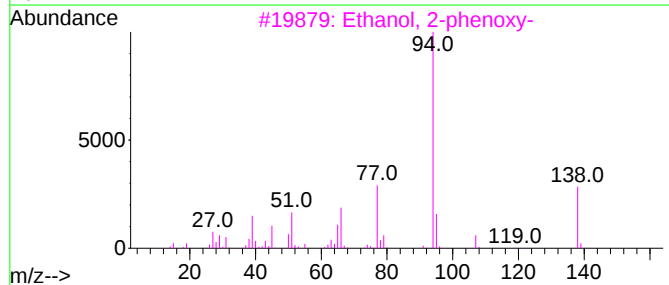
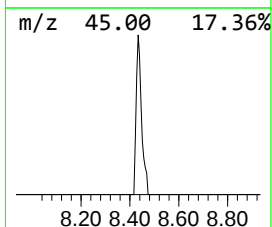
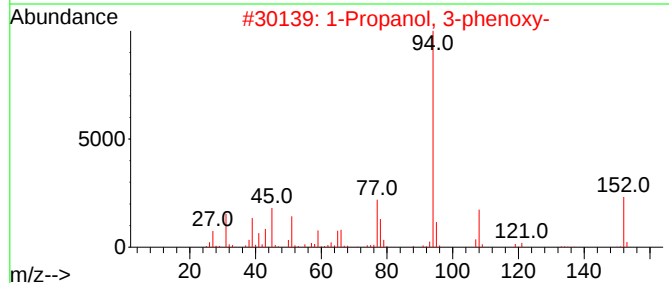
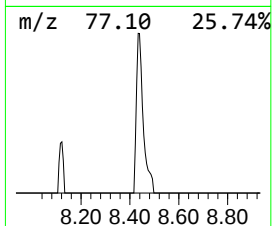
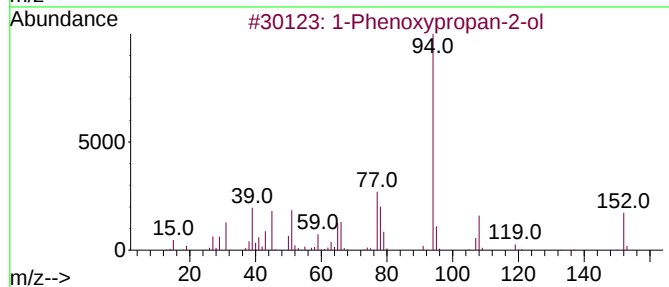
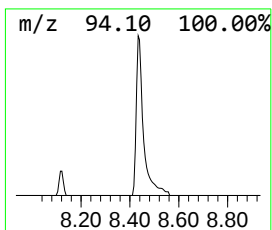
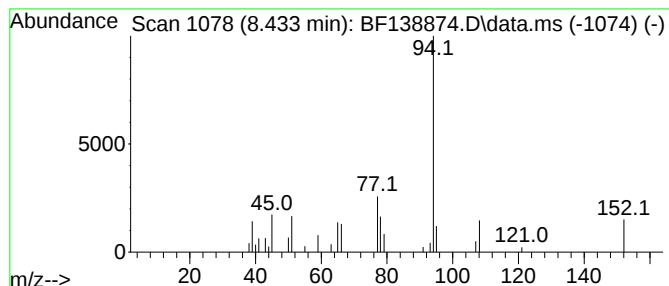
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Peak Number 2 1-Phenoxypropan-2-ol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.433	2.31 ng	42850	Naphthalene-d8	8.122

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	97
2		1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	91
3		Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	64
4		1-Propanol, 2-phenoxy-	152	C9H12O2	004169-04-4	64
5		Benzene, (1,1-dimethylethoxy)-	150	C10H14O	006669-13-2	53



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.051	4.8	ng	70897	1	6.839	297978	20.0
1-Phenoxypropan...	8.433	2.3	ng	42850	2	8.122	370399	20.0