

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111021\
 Data File : BF126099.D
 Acq On : 10 Nov 2021 10:15
 Operator : JU/CG
 Sample : PB140520BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB140520BL

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-BF110321.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF126099.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.205	12	20	26	rVB	184663	259531	2.61%	0.477%
2	5.481	571	577	580	rBV	6396915	6790073	68.18%	12.491%
3	6.487	741	748	751	rBV	6036573	6863625	68.92%	12.626%
4	6.851	806	810	814	rBV	1809132	1420228	14.26%	2.613%
5	7.416	900	906	909	rBV	4307794	4559297	45.78%	8.387%
6	8.134	1023	1028	1032	rBV	2042174	1812095	18.20%	3.333%
7	9.216	1205	1212	1214	rBV	8697574	8782922	88.20%	16.157%
8	9.886	1321	1326	1330	rBV	2460319	2217264	22.27%	4.079%
9	10.680	1455	1461	1464	rBV	6264815	6018388	60.43%	11.071%
10	11.375	1575	1579	1582	rBV	2627938	2298772	23.08%	4.229%
11	12.963	1844	1849	1852	rBV	10511190	9958465	100.00%	18.319%
12	14.010	2023	2027	2037	rBV	2131670	1928903	19.37%	3.548%
13	15.474	2270	2276	2281	rBV	1117500	1450836	14.57%	2.669%

Sum of corrected areas: 54360399

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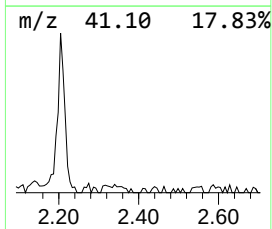
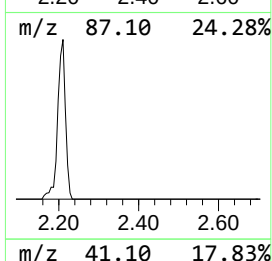
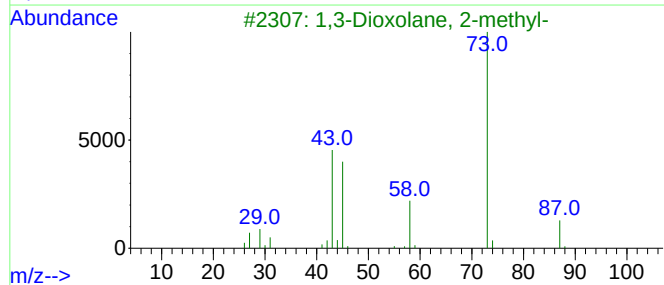
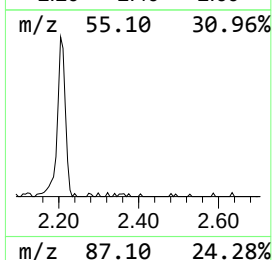
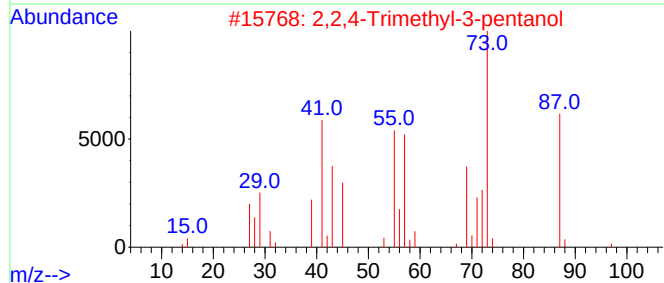
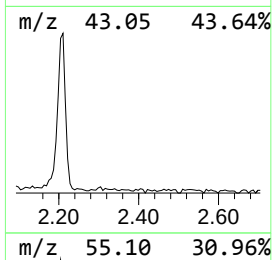
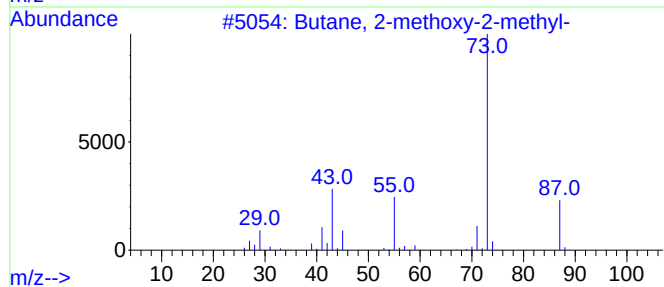
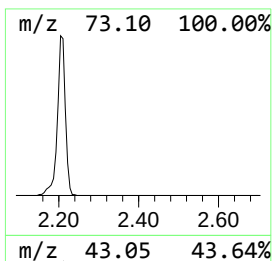
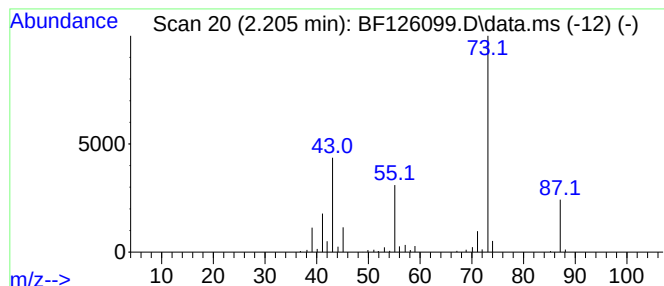
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.205	3.65 ng	259531	1,4-Dichlorobenzene-d4	6.851

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	37
4			3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	25
5			3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	17



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	2.205	3.6	ng	259531	1	6.851	1420230	20.0