

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011822\
 Data File : BP008838.D
 Acq On : 18 Jan 2022 14:06
 Operator : CG/JU
 Sample : PB142091BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB142091BL

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_P\Methods\8270E-BP011422.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP008838.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.058	8	12	18	rVB	225559	270765	1.56%	0.342%
2	5.440	410	417	440	rBV	5340848	8001747	46.16%	10.098%
3	7.034	680	688	705	rBV	4634396	7926242	45.73%	10.003%
4	7.852	819	827	843	rBV	1105039	1822932	10.52%	2.300%
5	9.010	1016	1024	1045	rBV	2935093	5204141	30.02%	6.567%
6	10.652	1295	1303	1323	rBV	1318070	2457579	14.18%	3.101%
7	13.116	1714	1722	1745	rBV	8839080	13351520	77.02%	16.849%
8	14.492	1948	1956	1976	rBV	2116397	3367803	19.43%	4.250%
9	15.987	2203	2210	2245	rBV	5894207	9442867	54.47%	11.917%
10	17.245	2418	2424	2438	rBV	2526120	3839906	22.15%	4.846%
11	19.810	2854	2860	2869	rBV	14171537	17334375	100.00%	21.875%
12	21.339	3115	3120	3131	rBV	2294915	3292966	19.00%	4.156%
13	23.763	3526	3532	3548	rVB2	1331082	2928664	16.90%	3.696%

Sum of corrected areas: 79241507

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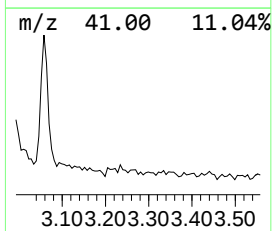
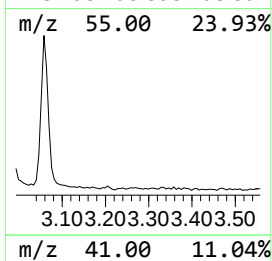
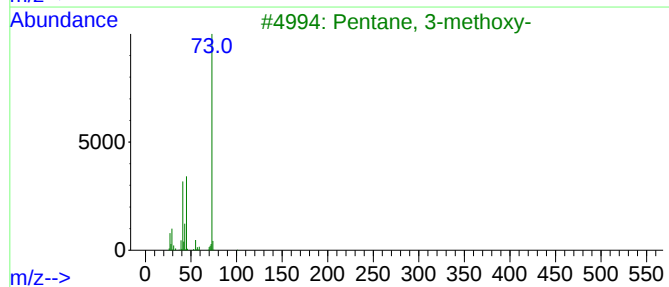
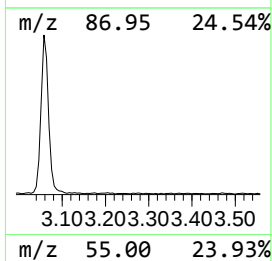
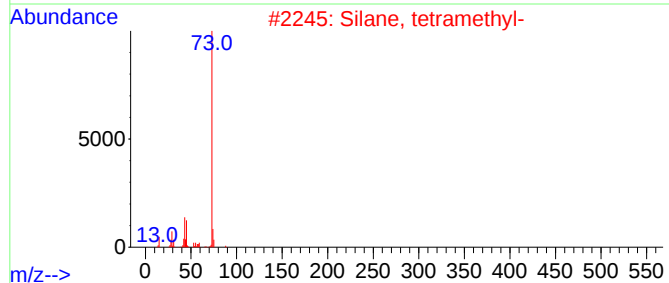
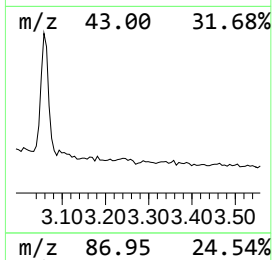
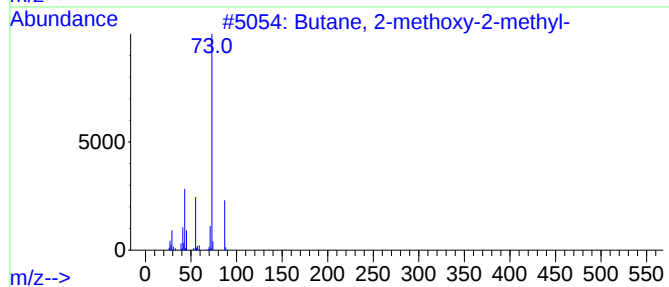
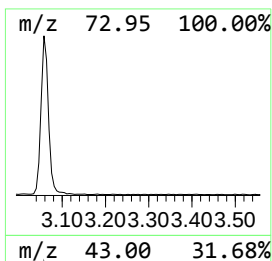
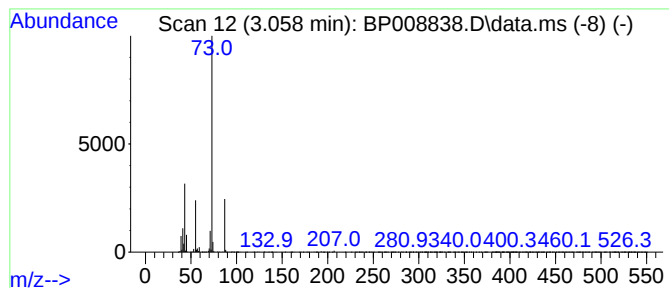
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
3.058	2.97 ng	270765	1,4-Dichlorobenzene-d4	7.852

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Silane, tetramethyl-	88	C4H12Si	000075-76-3	25
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.058	3.0	ng	270765	1	7.852	1822930	20.0