

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032922\
 Data File : BM034441.D
 Acq On : 30 Mar 2022 15:23
 Operator : CG/JU
 Sample : PB143666BL
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK666

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 03/31/2022
 Supervised By :Yogesh Patel 04/02/2022

Quant Time: Mar 31 02:21:05 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM032822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 28 18:27:40 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.925	152	3478	0.400	ng/u1	-0.01
4) Naphthalene-d8	10.735	136	9238	0.400	ng/u1	#-0.02
9) Acenaphthene-d10	14.569	164	5153	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.308	188	9429	0.400	ng/u1	-0.02
17) Chrysene-d12	21.507	240	5951	0.400	ng/u1	# 0.00
23) Perylene-d12	23.895	264	19400m	0.400	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.301	96	32109	6.206	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.341	152	4931	0.396	ng/u1	0.00
18) Fluoranthene-d10	19.339	212	8923	0.481	ng/u1	0.00

Target Compounds	Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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