

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032922\
 Data File : BM034456.D
 Acq On : 31 Mar 2022 01:00
 Operator : CG/JU
 Sample : PB143670BL
 Misc :
 ALS Vial : 61 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK670

Manual Integrations
 APPROVED

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

Quant Time: Mar 31 09:31:15 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	2976	0.400	ng/u1	-0.01
4) Naphthalene-d8	10.730	136	7972	0.400	ng/u1	#-0.02
9) Acenaphthene-d10	14.564	164	4276	0.400	ng/u1	-0.01
13) Phenanthrene-d10	17.308	188	8182	0.400	ng/u1	-0.02
17) Chrysene-d12	21.524	240	5206	0.400	ng/u1	# 0.02
23) Perylene-d12	23.886	264	18015m	0.400	ng/u1	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.301	96	22736	5.136	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.357	152	3220	0.300	ng/u1	0.02
18) Fluoranthene-d10	19.334	212	6624	0.408	ng/u1	-0.01

Target Compounds	Qvalue

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

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