

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062022\
 Data File : BM035579.D
 Acq On : 21 Jun 2022 19:48
 Operator : CG/JU
 Sample : PB145592BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK592

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 06/22/2022
 Supervised By :Yogesh Patel 06/25/2022

Quant Time: Jun 22 01:33:12 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM061822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 20 01:56:13 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.825	152	5725	0.400	ng/u1	-0.03
4) Naphthalene-d8	10.611	136	19741	0.400	ng/u1	-0.03
9) Acenaphthene-d10	14.455	164	11843	0.400	ng/u1	-0.01
13) Phenanthrene-d10	17.204	188	25301m	0.400	ng/u1	-0.03
17) Chrysene-d12	21.415	240	16392	0.400	ng/u1	0.00
23) Perylene-d12	23.733	264	18162	0.400	ng/u1	#-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	38163	4.738	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.250	152	8181	0.279	ng/u1	0.00
18) Fluoranthene-d10	19.238	212	17589	0.319	ng/u1	-0.01

Target Compounds	Qvalue

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

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