

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031722\
 Data File : BP009449.D
 Acq On : 17 Mar 2022 18:17
 Operator : CG/JU
 Sample : PB143392BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB143392BL

Quant Time: Mar 17 18:48:39 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 17 18:26:53 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.799	152	272013	20.000	ng	0.00
21) Naphthalene-d8	10.593	136	1056366	20.000	ng	0.00
39) Acenaphthene-d10	14.445	164	678474	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	1354560	20.000	ng	0.00
76) Chrysene-d12	21.316	240	1227764	20.000	ng	0.00
86) Perylene-d12	23.721	264	1197094	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.405	112	2037119	130.754	ng	0.00
7) Phenol-d6	6.993	99	2520898	119.647	ng	0.00
23) Nitrobenzene-d5	8.957	82	1640739	82.537	ng	0.00
42) 2,4,6-Tribromophenol	15.945	330	1247119	111.390	ng	0.00
45) 2-Fluorobiphenyl	13.063	172	4025132	82.243	ng	0.00
79) Terphenyl-d14	19.780	244	6013231	87.909	ng	0.00

Target Compounds	Qvalue

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

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