

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032922\
 Data File : BP009590.D
 Acq On : 29 Mar 2022 12:00
 Operator : CG/JU
 Sample : PB143645BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB143645BL

Quant Time: Mar 30 04:07:32 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 30 04:00:13 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	244025	20.000	ng	0.00
21) Naphthalene-d8	10.563	136	925264	20.000	ng	0.00
39) Acenaphthene-d10	14.416	164	554292	20.000	ng	0.00
64) Phenanthrene-d10	17.180	188	1132866	20.000	ng	0.00
76) Chrysene-d12	21.298	240	1106393	20.000	ng	0.00
86) Perylene-d12	23.692	264	1237602	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.369	112	2122621	151.868	ng	0.00
7) Phenol-d6	6.957	99	2580847	136.541	ng	0.00
23) Nitrobenzene-d5	8.922	82	1610983	92.523	ng	0.00
42) 2,4,6-Tribromophenol	15.916	330	1035821	113.244	ng	0.00
45) 2-Fluorobiphenyl	13.034	172	3661527	91.575	ng	0.00
79) Terphenyl-d14	19.763	244	5484917	88.981	ng	0.00

Target Compounds	Qvalue

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

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