

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031822\
 Data File : BP009457.D
 Acq On : 18 Mar 2022 14:54
 Operator : CG/JU
 Sample : PB143171BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB143171BL

Quant Time: Mar 18 17:35:15 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 18 17:04:24 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.799	152	310285	20.000	ng	0.00
21) Naphthalene-d8	10.593	136	1178146	20.000	ng	0.00
39) Acenaphthene-d10	14.439	164	758411	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	1531765	20.000	ng	0.00
76) Chrysene-d12	21.316	240	1316762	20.000	ng	0.00
86) Perylene-d12	23.715	264	1270003	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.405	112	2314801	130.251	ng	0.00
7) Phenol-d6	6.987	99	2847911	118.495	ng	0.00
23) Nitrobenzene-d5	8.952	82	1841303	83.052	ng	0.00
42) 2,4,6-Tribromophenol	15.939	330	1381415	110.380	ng	0.00
45) 2-Fluorobiphenyl	13.063	172	4424116	80.868	ng	0.00
79) Terphenyl-d14	19.780	244	6433034	87.689	ng	0.00

Target Compounds	Qvalue

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

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