

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011222\
 Data File : BP008771.D
 Acq On : 12 Jan 2022 13:23
 Operator : CG/JU
 Sample : PB141938BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB141938BL

Quant Time: Jan 12 16:21:31 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP010622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 12 00:18:56 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	339442	20.000	ng	0.00
21) Naphthalene-d8	10.669	136	1283997	20.000	ng	0.00
39) Acenaphthene-d10	14.504	164	848239	20.000	ng	0.00
64) Phenanthrene-d10	17.257	188	1877935	20.000	ng	0.00
76) Chrysene-d12	21.345	240	2035180	20.000	ng	0.00
86) Perylene-d12	23.774	264	2161399	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.452	112	2570124	118.569	ng	0.00
7) Phenol-d6	7.046	99	3094041	106.328	ng	0.00
23) Nitrobenzene-d5	9.028	82	1871917	82.564	ng	0.00
42) 2,4,6-Tribromophenol	15.998	330	1528702	106.822	ng	0.00
45) 2-Fluorobiphenyl	13.128	172	4920415	81.552	ng	0.00
79) Terphenyl-d14	19.822	244	8995204	90.070	ng	0.00

Target Compounds	Qvalue

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

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