

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF041422\
 Data File : BF127793.D
 Acq On : 14 Apr 2022 14:39
 Operator : CG\JU
 Sample : PB144112BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB144112BL

Quant Time: Apr 15 05:05:58 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF040622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 13 16:40:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.957	152	168545	20.000	ng	0.00
21) Naphthalene-d8	8.239	136	626846	20.000	ng	# 0.00
39) Acenaphthene-d10	9.998	164	358934	20.000	ng	0.00
64) Phenanthrene-d10	11.492	188	670685	20.000	ng	0.00
76) Chrysene-d12	14.139	240	557539	20.000	ng	0.00
86) Perylene-d12	15.657	264	442051	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.593	112	1244291	119.959	ng	0.01
7) Phenol-d6	6.581	99	1600348	117.936	ng	0.00
23) Nitrobenzene-d5	7.522	82	1147705	81.625	ng	0.00
42) 2,4,6-Tribromophenol	10.792	330	492339	128.898	ng	0.00
45) 2-Fluorobiphenyl	9.316	172	1978419	77.975	ng	0.00
79) Terphenyl-d14	13.086	244	2357804	69.803	ng	0.00

Target Compounds	Qvalue

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

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