

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122921\
 Data File : BP008592.D
 Acq On : 29 Dec 2021 10:14
 Operator : CG/JU
 Sample : PB141666BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB141666BL

Quant Time: Dec 29 13:37:53 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP122821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 29 13:29:13 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.899	152	688132	20.000	ng	0.00
21) Naphthalene-d8	10.699	136	2776856	20.000	ng	0.00
39) Acenaphthene-d10	14.528	164	2014561	20.000	ng	0.00
64) Phenanthrene-d10	17.275	188	4490998	20.000	ng	0.00
76) Chrysene-d12	21.357	240	3888019	20.000	ng	0.00
86) Perylene-d12	23.780	264	3398516	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.475	112	4832027	128.169	ng	0.00
7) Phenol-d6	7.063	99	6108063	116.399	ng	0.00
23) Nitrobenzene-d5	9.052	82	3630888	77.785	ng	0.00
42) 2,4,6-Tribromophenol	16.016	330	3404401	124.904	ng	0.00
45) 2-Fluorobiphenyl	13.157	172	10816897	81.113	ng	0.00
79) Terphenyl-d14	19.827	244	13966213	83.063	ng	0.00

Target Compounds	Qvalue

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

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