

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110321\
 Data File : BP007822.D
 Acq On : 03 Nov 2021 14:44
 Operator : CG/JU
 Sample : PB140419BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB140419BL

Quant Time: Nov 03 15:18:47 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 02 19:18:55 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	143037	20.00	ng	0.00
21) Naphthalene-d8	10.722	136	598407	20.00	ng	0.00
39) Acenaphthene-d10	14.557	164	439782	20.00	ng	0.00
64) Phenanthrene-d10	17.316	188	1070741	20.00	ng	0.00
76) Chrysene-d12	21.404	240	1339848	20.00	ng	-0.01
86) Perylene-d12	23.839	264	1511894	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.499	112	1309619	155.28	ng	0.00
7) Phenol-d6	7.093	99	1678180	137.47	ng	0.00
23) Nitrobenzene-d5	9.081	82	985864	88.50	ng	0.00
42) 2,4,6-Tribromophenol	16.051	330	864006	137.66	ng	0.00
45) 2-Fluorobiphenyl	13.187	172	2558925	84.74	ng	0.00
79) Terphenyl-d14	19.881	244	5420901	82.39	ng	0.00

Target Compounds	Qvalue

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

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