

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070921\
 Data File : BP006159.D
 Acq On : 09 Jul 2021 14:02
 Operator : CG/JU
 Sample : PB137592BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB137592BL

Quant Time: Jul 09 15:29:24 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP070721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 07 16:08:13 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	150586	20.000	ng	0.00
21) Naphthalene-d8	10.681	136	537946	20.000	ng	0.00
39) Acenaphthene-d10	14.516	164	342317	20.000	ng	0.00
64) Phenanthrene-d10	17.269	188	680104	20.000	ng	# 0.00
76) Chrysene-d12	21.345	240	712852	20.000	ng	# 0.00
86) Perylene-d12	23.751	264	880135	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	1355784	154.916	ng	-0.01
7) Phenol-d6	7.058	99	1544588	133.428	ng	0.00
23) Nitrobenzene-d5	9.040	82	875190	102.764	ng	0.00
42) 2,4,6-Tribromophenol	16.010	330	647012	158.756	ng	0.00
45) 2-Fluorobiphenyl	13.140	172	1949151	84.177	ng	0.00
79) Terphenyl-d14	19.816	244	3494172	96.643	ng	0.00

Target Compounds	Qvalue

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

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