

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091621\
 Data File : BP006996.D
 Acq On : 16 Sep 2021 15:26
 Operator : CG/JU
 Sample : PB139068BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB139068BL

Quant Time: Sep 16 16:06:44 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 16 15:16:52 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.916	152	336016	20.000	ng	0.00
21) Naphthalene-d8	10.716	136	1304407	20.000	ng	# 0.00
39) Acenaphthene-d10	14.545	164	811953	20.000	ng	0.00
64) Phenanthrene-d10	17.292	188	1780513	20.000	ng	# 0.00
76) Chrysene-d12	21.374	240	1779546	20.000	ng	# 0.00
86) Perylene-d12	23.792	264	1789271	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	2668909	129.700	ng	0.00
7) Phenol-d6	7.075	99	3230473	115.019	ng	0.00
23) Nitrobenzene-d5	9.075	82	1805996	77.762	ng	0.00
42) 2,4,6-Tribromophenol	16.034	330	1252185	128.236	ng	0.00
45) 2-Fluorobiphenyl	13.175	172	4543082	77.243	ng	0.00
79) Terphenyl-d14	19.851	244	7751259	83.570	ng	0.00

Target Compounds	Qvalue

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

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