

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091621\
 Data File : BP006999.D
 Acq On : 16 Sep 2021 17:08
 Operator : CG/JU
 Sample : PB139068BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB139068BL

Quant Time: Sep 16 17:43:23 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	7.922	152	374512	20.000	ng	0.00
21) Naphthalene-d8	10.722	136	1440468	20.000	ng	# 0.00
39) Acenaphthene-d10	14.545	164	876107	20.000	ng	0.00
64) Phenanthrene-d10	17.292	188	1821554	20.000	ng	# 0.00
76) Chrysene-d12	21.368	240	1796824	20.000	ng	#-0.01
86) Perylene-d12	23.786	264	1822405	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.493	112	2951174	128.675	ng	0.00
7) Phenol-d6	7.081	99	3576103	114.237	ng	0.00
23) Nitrobenzene-d5	9.081	82	2014223	78.536	ng	0.00
42) 2,4,6-Tribromophenol	16.033	330	1269422	120.482	ng	0.00
45) 2-Fluorobiphenyl	13.175	172	4921375	77.547	ng	0.00
79) Terphenyl-d14	19.851	244	7759773	82.858	ng	0.00

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

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