

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092421\
 Data File : BP007158.D
 Acq On : 24 Sep 2021 14:36
 Operator : CG/JU
 Sample : M3921-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SPOILS-REPORT-1

Quant Time: Sep 24 17:06:09 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 23 11:17:25 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	228858	20.000	ng	0.00
21) Naphthalene-d8	10.687	136	872097	20.000	ng	# 0.00
39) Acenaphthene-d10	14.522	164	562630	20.000	ng	0.00
64) Phenanthrene-d10	17.269	188	1219914	20.000	ng	# 0.00
76) Chrysene-d12	21.357	240	1252994	20.000	ng	# 0.00
86) Perylene-d12	23.768	264	1309627	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.463	112	1306480	95.560	ng	0.00
7) Phenol-d6	7.058	99	1617815	86.579	ng	0.00
23) Nitrobenzene-d5	9.046	82	922534	61.604	ng	-0.01
42) 2,4,6-Tribromophenol	16.010	330	759434	104.115	ng	0.00
45) 2-Fluorobiphenyl	13.145	172	2394746	60.977	ng	0.00
79) Terphenyl-d14	19.827	244	3857920	59.001	ng	0.00

Target Compounds	Qvalue

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

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