

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080421\
 Data File : BP006489.D
 Acq On : 04 Aug 2021 11:06
 Operator : CG/JU
 Sample : PB138117BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB138117BL

Quant Time: Aug 04 11:45:56 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 03 15:41:30 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.757	152	190952	20.000	ng	0.00
21) Naphthalene-d8	10.540	136	809740	20.000	ng	# 0.00
39) Acenaphthene-d10	14.387	164	446763	20.000	ng	0.00
64) Phenanthrene-d10	17.139	188	876362	20.000	ng	# 0.00
76) Chrysene-d12	21.233	240	781612	20.000	ng	# 0.00
86) Perylene-d12	23.563	264	777795	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.358	112	1828036	147.040	ng	0.00
7) Phenol-d6	6.934	99	2352085	133.185	ng	0.00
23) Nitrobenzene-d5	8.916	82	1539265	87.742	ng	0.00
42) 2,4,6-Tribromophenol	15.881	330	573151	122.752	ng	0.00
45) 2-Fluorobiphenyl	13.010	172	2648959	88.709	ng	0.00
79) Terphenyl-d14	19.716	244	3602509	92.356	ng	0.00

Target Compounds	Qvalue

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

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