

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111821\
 Data File : BP008008.D
 Acq On : 18 Nov 2021 10:03
 Operator : CG/JU
 Sample : PB140849BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB140849BL

Quant Time: Nov 18 10:51:59 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 15 05:36:49 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.834	152	115944	20.000	ng	0.00
21) Naphthalene-d8	10.634	136	463545	20.000	ng	0.00
39) Acenaphthene-d10	14.475	164	320650	20.000	ng	0.00
64) Phenanthrene-d10	17.233	188	740697	20.000	ng	# 0.00
76) Chrysene-d12	21.322	240	925665	20.000	ng	#-0.01
86) Perylene-d12	23.721	264	1071260	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	1039197	145.399	ng	0.00
7) Phenol-d6	7.016	99	1277899	134.399	ng	0.00
23) Nitrobenzene-d5	8.999	82	727827	84.023	ng	0.00
42) 2,4,6-Tribromophenol	15.969	330	689611	136.691	ng	0.00
45) 2-Fluorobiphenyl	13.098	172	1863952	81.541	ng	0.00
79) Terphenyl-d14	19.798	244	3721798	81.528	ng	0.00

Target Compounds	Qvalue

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

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