

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021422\
 Data File : BP009132.D
 Acq On : 14 Feb 2022 13:34
 Operator : CG/JU
 Sample : PB142649BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB142649BL

Quant Time: Feb 14 15:08:31 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 07 15:13:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	222957	20.000	ng	-0.01
21) Naphthalene-d8	10.593	136	848053	20.000	ng	-0.01
39) Acenaphthene-d10	14.440	164	597762	20.000	ng	0.00
64) Phenanthrene-d10	17.198	188	1278168	20.000	ng	0.00
76) Chrysene-d12	21.298	240	1101653	20.000	ng	0.00
86) Perylene-d12	23.698	264	959260	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	1871324	149.718	ng	0.00
7) Phenol-d6	6.993	99	2295029	139.351	ng	0.00
23) Nitrobenzene-d5	8.958	82	1454908	96.748	ng	-0.01
42) 2,4,6-Tribromophenol	15.945	330	1270481	159.184	ng	0.00
45) 2-Fluorobiphenyl	13.057	172	3969112	92.952	ng	-0.01
79) Terphenyl-d14	19.763	244	6114959	99.820	ng	0.00

Target Compounds	Qvalue

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

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