

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020122\
 Data File : BP008980.D
 Acq On : 01 Feb 2022 11:21
 Operator : CG/JU
 Sample : PB142379BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB142379BL

Quant Time: Feb 01 14:05:08 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 01 04:16:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.828	152	229108	20.000	ng	0.00
21) Naphthalene-d8	10.628	136	837651	20.000	ng	0.00
39) Acenaphthene-d10	14.469	164	467506	20.000	ng	0.00
64) Phenanthrene-d10	17.222	188	872813	20.000	ng	-0.01
76) Chrysene-d12	21.316	240	893510	20.000	ng	-0.01
86) Perylene-d12	23.721	264	1073175	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	1744413	117.743	ng	0.00
7) Phenol-d6	7.010	99	2022871	112.754	ng	0.00
23) Nitrobenzene-d5	8.987	82	1290211	87.447	ng	-0.01
42) 2,4,6-Tribromophenol	15.963	330	765449	112.483	ng	-0.01
45) 2-Fluorobiphenyl	13.092	172	3077292	86.804	ng	0.00
79) Terphenyl-d14	19.786	244	3966309	74.928	ng	-0.01

Target Compounds	Qvalue

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

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