

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020222\
 Data File : BP009004.D
 Acq On : 02 Feb 2022 13:48
 Operator : CG/JU
 Sample : PB142428BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB142428BL

Quant Time: Feb 03 06:23:22 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 03 06:20:22 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.828	152	241630	20.000	ng	0.00
21) Naphthalene-d8	10.628	136	974479	20.000	ng	0.00
39) Acenaphthene-d10	14.469	164	661951	20.000	ng	0.00
64) Phenanthrene-d10	17.228	188	1425884	20.000	ng	0.00
76) Chrysene-d12	21.316	240	1340993	20.000	ng	0.00
86) Perylene-d12	23.727	264	1246686	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.417	112	1840166	117.769	ng	0.00
7) Phenol-d6	7.011	99	2267848	119.858	ng	0.00
23) Nitrobenzene-d5	8.987	82	1478365	86.130	ng	0.00
42) 2,4,6-Tribromophenol	15.969	330	1224587	127.093	ng	0.00
45) 2-Fluorobiphenyl	13.093	172	3995414	79.597	ng	0.00
79) Terphenyl-d14	19.786	244	6575158	82.763	ng	0.00

Target Compounds	Qvalue

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

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