

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020322\
 Data File : BP009022.D
 Acq On : 03 Feb 2022 11:09
 Operator : CG/JU
 Sample : PB142438BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB142438BL

Quant Time: Feb 03 13:30:01 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	7.822	152	240700	20.000	ng	0.00
21) Naphthalene-d8	10.622	136	865817	20.000	ng	0.00
39) Acenaphthene-d10	14.469	164	509841	20.000	ng	0.00
64) Phenanthrene-d10	17.227	188	924245	20.000	ng	0.00
76) Chrysene-d12	21.316	240	836650	20.000	ng	0.00
86) Perylene-d12	23.727	264	943744	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	1786905	114.803	ng	0.00
7) Phenol-d6	7.010	99	2098229	111.322	ng	0.00
23) Nitrobenzene-d5	8.987	82	1323215	86.766	ng	0.00
42) 2,4,6-Tribromophenol	15.969	330	794329	107.034	ng	0.00
45) 2-Fluorobiphenyl	13.087	172	3205607	82.915	ng	0.00
79) Terphenyl-d14	19.786	244	3727110	75.194	ng	0.00

Target Compounds	Qvalue

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

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