

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030822\
 Data File : BP009301.D
 Acq On : 08 Mar 2022 13:03
 Operator : CG/JU
 Sample : PB143112BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB143112BL

Quant Time: Mar 09 03:00:28 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 05 01:55:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.816	152	513392	20.000	ng	0.00
21) Naphthalene-d8	10.610	136	2009956	20.000	ng	0.00
39) Acenaphthene-d10	14.457	164	1181496	20.000	ng	0.00
64) Phenanthrene-d10	17.216	188	2199308	20.000	ng	0.00
76) Chrysene-d12	21.321	240	1780819	20.000	ng	0.00
86) Perylene-d12	23.721	264	1840848	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	4284095	121.806	ng	0.00
7) Phenol-d6	6.993	99	5332498	119.307	ng	0.00
23) Nitrobenzene-d5	8.969	82	3324758	91.108	ng	0.00
42) 2,4,6-Tribromophenol	15.951	330	1895747	132.880	ng	0.00
45) 2-Fluorobiphenyl	13.081	172	7407945	86.225	ng	0.00
79) Terphenyl-d14	19.792	244	8902285	94.337	ng	0.00

Target Compounds	Qvalue

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

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