

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030522\
 Data File : BP009277.D
 Acq On : 04 Mar 2022 20:58
 Operator : CG/JU
 Sample : PB143102BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB143102BL

Quant Time: Mar 05 02:04:17 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	7.822	152	389127	20.000	ng	0.00
21) Naphthalene-d8	10.616	136	1537094	20.000	ng	0.00
39) Acenaphthene-d10	14.457	164	929787	20.000	ng	0.00
64) Phenanthrene-d10	17.216	188	1962631	20.000	ng	0.00
76) Chrysene-d12	21.327	240	2027093	20.000	ng	0.00
86) Perylene-d12	23.733	264	2088156	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	4069085	152.638	ng	0.00
7) Phenol-d6	6.993	99	5045918	148.947	ng	0.00
23) Nitrobenzene-d5	8.975	82	3109181	111.411	ng	0.00
42) 2,4,6-Tribromophenol	15.951	330	1839193	163.816	ng	0.00
45) 2-Fluorobiphenyl	13.087	172	7064505	104.488	ng	0.00
79) Terphenyl-d14	19.798	244	11839576	110.221	ng	0.00

Target Compounds	Qvalue

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

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