

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030922\
 Data File : BP009319.D
 Acq On : 09 Mar 2022 12:04
 Operator : CG/JU
 Sample : PB143171BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB143171BL

Quant Time: Mar 10 04:08:45 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	469893	20.000	ng	0.00
21) Naphthalene-d8	10.616	136	1854948	20.000	ng	0.00
39) Acenaphthene-d10	14.457	164	1133157	20.000	ng	0.00
64) Phenanthrene-d10	17.215	188	2234471	20.000	ng	0.00
76) Chrysene-d12	21.321	240	1970204	20.000	ng	0.00
86) Perylene-d12	23.727	264	2031242	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	3817013	118.572	ng	0.00
7) Phenol-d6	6.992	99	4709715	115.128	ng	0.00
23) Nitrobenzene-d5	8.975	82	2987951	88.721	ng	0.00
42) 2,4,6-Tribromophenol	15.951	330	1922880	140.531	ng	0.00
45) 2-Fluorobiphenyl	13.086	172	6950220	84.348	ng	0.00
79) Terphenyl-d14	19.792	244	9306933	89.145	ng	0.00

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

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