

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040122\
 Data File : BP009653.D
 Acq On : 01 Apr 2022 12:14
 Operator : CG/JU
 Sample : PB143765BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB143765BL

Quant Time: Apr 01 22:42:48 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 31 04:36:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.728	152	264659	20.000	ng	-0.01
21) Naphthalene-d8	10.522	136	954764	20.000	ng	-0.01
39) Acenaphthene-d10	14.381	164	581480	20.000	ng	-0.01
64) Phenanthrene-d10	17.151	188	1201747	20.000	ng	0.00
76) Chrysene-d12	21.269	240	1171116	20.000	ng	0.00
86) Perylene-d12	23.645	264	1106111	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.340	112	2327701	153.556	ng	0.00
7) Phenol-d6	6.928	99	2842204	138.645	ng	0.00
23) Nitrobenzene-d5	8.893	82	1800814	100.230	ng	0.00
42) 2,4,6-Tribromophenol	15.887	330	1114833	116.184	ng	0.00
45) 2-Fluorobiphenyl	12.999	172	3842023	91.597	ng	-0.01
79) Terphenyl-d14	19.733	244	5832764	89.395	ng	0.00

Target Compounds	Qvalue

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

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