

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033122\
 Data File : BP009634.D
 Acq On : 31 Mar 2022 12:02
 Operator : CG/JU
 Sample : PB143709BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB143709BL

Quant Time: Mar 31 14:17:45 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	7.740	152	313017	20.000	ng	0.00
21) Naphthalene-d8	10.534	136	1317206	20.000	ng	0.00
39) Acenaphthene-d10	14.392	164	924304	20.000	ng	0.00
64) Phenanthrene-d10	17.157	188	2058906	20.000	ng	0.00
76) Chrysene-d12	21.274	240	1843155	20.000	ng	0.00
86) Perylene-d12	23.657	264	1703393	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.346	112	2624923	146.412	ng	0.00
7) Phenol-d6	6.934	99	3399434	140.209	ng	0.00
23) Nitrobenzene-d5	8.899	82	2231734	90.036	ng	0.00
42) 2,4,6-Tribromophenol	15.898	330	1949056	127.785	ng	0.00
45) 2-Fluorobiphenyl	13.010	172	5564248	83.454	ng	0.00
79) Terphenyl-d14	19.739	244	9637969	93.856	ng	0.00

Target Compounds	Qvalue

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

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