

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032222\
 Data File : BM034306.D
 Acq On : 22 Mar 2022 12:10
 Operator : CG/JU
 Sample : PB143452BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB143452BL

Quant Time: Mar 22 14:37:15 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM031522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 16 02:38:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.942	152	155637	20.000	ng	0.00
21) Naphthalene-d8	10.754	136	625840	20.000	ng	-0.01
39) Acenaphthene-d10	14.583	164	406938	20.000	ng	0.00
64) Phenanthrene-d10	17.318	188	845098	20.000	ng	-0.01
76) Chrysene-d12	21.489	240	844789	20.000	ng	-0.01
86) Perylene-d12	23.900	264	865058	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.496	112	1394630	160.753	ng	0.00
7) Phenol-d6	7.101	99	1822808	146.747	ng	0.00
23) Nitrobenzene-d5	9.101	82	1167614	90.834	ng	-0.01
42) 2,4,6-Tribromophenol	16.066	330	750183	136.770	ng	0.00
45) 2-Fluorobiphenyl	13.213	172	2673321	89.770	ng	0.00
79) Terphenyl-d14	19.942	244	4527832	93.131	ng	0.00

Target Compounds	Qvalue

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

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