

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110922\
 Data File : BM037434.D
 Acq On : 09 Nov 2022 12:11
 Operator : CG/JU
 Sample : PB148855BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB148855BL

Quant Time: Nov 10 05:46:54 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 09 02:08:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.816	152	352633	20.000	ng	0.00
21) Naphthalene-d8	10.616	136	1502512	20.000	ng	-0.01
39) Acenaphthene-d10	14.457	164	933133	20.000	ng	0.00
64) Phenanthrene-d10	17.198	188	1778933	20.000	ng	0.00
76) Chrysene-d12	21.374	240	1304985	20.000	ng	0.00
86) Perylene-d12	23.709	264	1167161	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	2723025	133.408	ng	0.00
7) Phenol-d6	6.998	99	3581948	129.301	ng	0.00
23) Nitrobenzene-d5	8.992	82	2180550	73.220	ng	0.00
42) 2,4,6-Tribromophenol	15.951	330	1602891	145.762	ng	0.00
45) 2-Fluorobiphenyl	13.080	172	5354264	76.240	ng	-0.01
79) Terphenyl-d14	19.821	244	6978685	95.899	ng	0.00

Target Compounds	Qvalue

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

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