

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071522\
 Data File : BM035842.D
 Acq On : 15 Jul 2022 22:28
 Operator : CG/JU
 Sample : N3652-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SED-3-(18-24)

Quant Time: Jul 15 23:48:50 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM071322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 14 01:21:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.904	152	345468	20.000	ng	0.00
21) Naphthalene-d8	10.710	136	1500248	20.000	ng	0.00
39) Acenaphthene-d10	14.539	164	880084	20.000	ng	0.00
64) Phenanthrene-d10	17.274	188	1668277	20.000	ng	0.00
76) Chrysene-d12	21.445	240	1189512	20.000	ng	-0.01
86) Perylene-d12	23.821	264	1106437	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.463	112	2438973	114.673	ng	0.00
7) Phenol-d6	7.069	99	3132082	117.729	ng	0.00
23) Nitrobenzene-d5	9.069	82	1840214	70.122	ng	0.00
42) 2,4,6-Tribromophenol	16.021	330	1038226	116.167	ng	0.00
45) 2-Fluorobiphenyl	13.169	172	3790190	60.061	ng	0.00
79) Terphenyl-d14	19.898	244	4201678	68.568	ng	0.00

Target Compounds	Qvalue

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

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