

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090222\
 Data File : BN021592.D
 Acq On : 03 Sep 2022 11:47
 Operator : CG/JU
 Sample : N4454-10
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MW-07-1.5-11.5-082922

Quant Time: Sep 05 02:38:46 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN082922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 02 00:53:31 2022
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 09/06/2022
 Supervised By :Jagrut Upadhyay 09/06/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.069	152	5230	0.400	ng	0.00
7) Naphthalene-d8	10.898	136	17670	0.400	ng	0.00
13) Acenaphthene-d10	14.718	164	10457	0.400	ng	0.00
19) Phenanthrene-d10	17.470	188	21950	0.400	ng	0.00
29) Chrysene-d12	21.655	240	16333	0.400	ng	# 0.00
35) Perylene-d12	24.173	264	12242	0.400	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.592	112	1593	0.156	ng	0.00
5) Phenol-d6	7.217	99	1253	0.096	ng	0.00
8) Nitrobenzene-d5	9.233	82	3707m	0.311	ng	-0.01
11) 2-Methylnaphthalene-d10	12.482	152	11105	0.279	ng	0.00
14) 2,4,6-Tribromophenol	16.208	330	1335	0.389	ng	-0.01
15) 2-Fluorobiphenyl	13.349	172	17522	0.383	ng	0.00
27) Fluoranthene-d10	19.497	212	20494	0.363	ng	0.00
31) Terphenyl-d14	20.088	244	17327	0.472	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.403	88	5577	0.853	ng	# 94
17) Acenaphthene	14.782	154	11950	0.346	ng	99
34) Bis(2-ethylhexyl)phtha...	21.557	149	4192	0.183	ng	99

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

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