

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071421\
 Data File : BN015569.D
 Acq On : 14 Jul 2021 18:27
 Operator : CG/JU
 Sample : M2953-02DL 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MW-11B-070621DL

Quant Time: Jul 15 01:05:10 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN063021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 14 11:20:16 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.817	152	16503	0.400	ng	# 0.00
7) Naphthalene-d8	10.596	136	77669	0.400	ng	# 0.00
13) Acenaphthene-d10	14.446	164	42900	0.400	ng	0.01
19) Phenanthrene-d10	17.188	188	83871	0.400	ng	0.00
29) Chrysene-d12	21.386	240	56679	0.400	ng	# 0.00
35) Perylene-d12	23.744	264	30200	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.429	112	981	0.023	ng	0.00
5) Phenol-d6	6.987	99	663	0.013	ng	0.00
8) Nitrobenzene-d5	8.962	82	3240	0.074	ng	0.00
11) 2-Methylnaphthalene-d10	12.194	152	8173	0.064	ng	0.00
14) 2,4,6-Tribromophenol	15.943	330	735	0.259	ng	0.01
15) 2-Fluorobiphenyl	13.063	172	11409	0.065	ng	0.00
27) Fluoranthene-d10	19.226	212	19315	0.078	ng	0.00
31) Terphenyl-d14	19.832	244	14440	0.109	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.382	88	41819	5.405	ng	# 62
23) Atrazine	16.652	200	89	0.175	ng	# 47
24) Pentachlorophenol	16.835	266	1245	0.366	ng	# 35
34) Bis(2-ethylhexyl)phtha...	21.301	149	6092	0.236	ng	# 87

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

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