

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN040422\
 Data File : BN019346.D
 Acq On : 05 Apr 2022 10:08
 Operator : CG/JU
 Sample : SP5858
 Misc : 8270 SIM SURROGATE
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SP5858

Quant Time: Apr 05 10:46:28 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN040422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 04 17:40:00 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.833	152	4492	0.400	ng	0.00
7) Naphthalene-d8	10.637	136	13409	0.400	ng	# 0.01
13) Acenaphthene-d10	14.463	164	8412	0.400	ng	0.00
19) Phenanthrene-d10	17.205	188	15958	0.400	ng	0.00
29) Chrysene-d12	21.395	240	12056	0.400	ng	# 0.00
35) Perylene-d12	23.738	264	9671	0.400	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.384	112	5082	0.422	ng	0.00
5) Phenol-d6	7.002	99	4488	0.310	ng	0.01
8) Nitrobenzene-d5	8.993	82	3358	0.276	ng	0.01
11) 2-Methylnaphthalene-d10	12.227	152	8310	0.353	ng	0.00
14) 2,4,6-Tribromophenol	15.964	330	1268	0.248	ng	0.01
15) 2-Fluorobiphenyl	13.095	172	11327	0.324	ng	0.00
27) Fluoranthene-d10	19.240	212	19325	0.376	ng	0.00
31) Terphenyl-d14	19.834	244	10359	0.395	ng	0.00

Target Compounds Qvalue

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

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