

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041921\
 Data File : BN014325.D
 Acq On : 19 Apr 2021 12:55
 Operator : CG/JU
 Sample : M2061-02DL 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GWBR3-225-250-041521DL

Quant Time: Apr 19 13:27:55 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN041621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 19 10:42:58 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.700	152	7706	0.40	ng	0.00
7) Naphthalene-d8	10.467	136	28182	0.40	ng	# 0.00
13) Acenaphthene-d10	14.321	164	14361	0.40	ng	0.00
19) Phenanthrene-d10	17.067	188	31818	0.40	ng	0.00
29) Chrysene-d12	21.269	240	26193	0.40	ng	# 0.00
36) Perylene-d12	23.500	264	22727	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.301	112	299	0.02	ng	0.00
5) Phenol-d6	6.863	99	213	0.01	ng	0.00
8) Nitrobenzene-d5	8.832	82	1008	0.06	ng	0.00
11) 2-Methylnaphthalene-d10	12.062	152	2174	0.05	ng	0.00
14) 2,4,6-Tribromophenol	15.818	330	314	0.06	ng	0.00
15) 2-Fluorobiphenyl	12.951	172	3378	0.06	ng	0.00
27) Fluoranthene-d10	19.106	212	5087	0.06	ng	0.00
31) Terphenyl-d14	19.712	244	5943	0.11	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.255	88	19581	2.00	ng	# 91
34) Bis(2-ethylhexyl)phtha...	21.194	149	2008	0.10	ng	97
35) Indeno(1,2,3-cd)pyrene	25.735	276	2093	0.02	ng	96
37) Benzo(b)fluoranthene	22.833	252	2265	0.02	ng	# 33
41) Benzo(g,h,i)perylene	26.427	276	2047	0.02	ng	# 77

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

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