

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN011322\
 Data File : BN018399.D
 Acq On : 13 Jan 2022 17:25
 Operator : CG/JU
 Sample : N1122-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GWBR3D-PDB-472-493-011222

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 01/14/2022
 Supervised By : Jagrut Upadhyay 01/14/2022

Quant Time: Jan 13 20:20:58 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN010722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 10 13:58:28 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.615	152	26174	0.400	ng	0.00
7) Naphthalene-d8	10.381	136	125924	0.400	ng	#-0.03
13) Acenaphthene-d10	14.243	164	68481	0.400	ng	-0.01
19) Phenanthrene-d10	16.981	188	115662	0.400	ng	#-0.01
29) Chrysene-d12	21.185	240	109926	0.400	ng	# 0.00
35) Perylene-d12	23.406	264	105351	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.236	112	10372	0.161	ng	0.00
5) Phenol-d6	6.812	99	6853	0.106	ng	0.00
8) Nitrobenzene-d5	8.759	82	31864	0.456	ng	-0.01
11) 2-Methylnaphthalene-d10	11.980	152	59876	0.304	ng	-0.02
14) 2,4,6-Tribromophenol	15.740	330	13287	0.551	ng	-0.01
15) 2-Fluorobiphenyl	12.873	172	85323	0.345	ng	-0.01
27) Fluoranthene-d10	19.023	212	130283	0.394	ng	0.00
31) Terphenyl-d14	19.629	244	105540	0.444	ng	-0.01
Target Compounds						
2) 1,4-Dioxane	3.161	88	93431m	6.346	ng	Qvalue
24) Pentachlorophenol	16.652	266	857	0.361	ng	96
25) Phenanthrene	17.030	178	11714	0.036	ng	97
34) Bis(2-ethylhexyl)phtha...	21.121	149	176780	0.818	ng	# 96
40) Dibenzo(a,h)anthracene	25.692	278	1318	0.043	ng	# 1

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

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