

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102918\
 Data File : BN003346.D
 Acq On : 29 Oct 2018 23:38
 Operator : MJ/SJ
 Sample : J5692-08
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SA-PZ143I-20181024

Quant Time: Oct 30 07:40:58 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN102618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 29 16:15:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	18801	0.40	ng	0.00
7) Naphthalene-d8	10.66	136	82767	0.40	ng	-0.01
13) Acenaphthene-d10	14.49	164	50363	0.40	ng	-0.01
19) Phenanthrene-d10	17.24	188	118795	0.40	ng	0.00
27) Chrysene-d12	21.42	240	127465	0.40	ng	0.00
34) Perylene-d12	23.73	264	109155	0.40	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.42	112	7918	0.17	ng	0.00
5) Phenol-d6	7.01	99	5915	0.10	ng	0.00
8) Nitrobenzene-d5	9.02	82	15507	0.43	ng	0.00
11) 2-Methylnaphthalene-d10	12.24	152	42512	0.32	ng	0.00
14) 2,4,6-Tribromophenol	15.98	330	8581	0.43	ng	0.00
15) 2-Fluorobiphenyl	13.12	172	71947	0.33	ng	-0.01
25) Fluoranthene-d10	19.26	212	106911	0.31	ng	0.00
29) Terphenyl-d14	19.86	244	113577	0.38	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.34	88	4974	0.204	ng	Qvalue # 75
32) Bis(2-ethylhexyl)phthalate	21.31	149	9139	0.092	ng	# 99
35) Benzo(b)fluoranthene	23.02	252	10380	0.030	ng	# 68
36) Benzo(k)fluoranthene	23.08	252	7304	0.022	ng	# 64
37) Benzo(a)pyrene	23.63	252	6558	0.021	ng	# 56
39) Benzo(g,h,i)perylene	26.81	276	6022	0.021	ng	# 89

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

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