

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101518\
 Data File : BF109891.D
 Acq On : 15 Oct 2018 20:47
 Operator : JU/SJ
 Sample : J5198-13
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-02-101218-B

Manual Integrations
 APPROVED

Sohil
 10/16/2018 9:57:36 AM

Quant Time: Oct 16 01:55:15 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 15 18:09:48 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	217785	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	793560	20.00	ng	0.00
39) Acenaphthene-d10	10.02	164	303101	20.00	ng	0.00
64) Phenanthrene-d10	11.51	188	494119	20.00	ng	0.00
76) Chrysene-d12	14.15	240	518386	20.00	ng	0.00
87) Perylene-d12	15.67	264	416789	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	1178441	85.77	ng	0.00
7) Phenol-d6	6.59	99	1435183	84.12	ng	0.00
23) Nitrobenzene-d5	7.53	82	827068	70.21	ng	-0.01
42) 2,4,6-Tribromophenol	10.80	330	218093	83.11	ng	0.00
45) 2-Fluorobiphenyl	9.33	172	1237231	65.13	ng	0.00
79) Terphenyl-d14	13.09	244	1019422	41.30	ng	0.00
Target Compounds						
10) Phenol	6.60	94	51731	2.808	ng	Qvalue # 59
50) Dimethylphthalate	9.72	163	141568	6.596	ng	100
71) Phenanthrene	11.53	178	169809	6.638	ng	98
75) Fluoranthene	12.72	202	136706	5.349	ng	98
78) Pyrene	12.95	202	175416	4.607	ng	99
81) Benzo(a)anthracene	14.14	228	68513m	2.244	ng	
83) Chrysene	14.17	228	83082m	2.859	ng	
88) Benzo(b)fluoranthene	15.22	252	59802m	2.489	ng	

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

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