

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100518\
 Data File : BF109609.D
 Acq On : 5 Oct 2018 13:49
 Operator : JU/SJ
 Sample : J5150-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSSI-0926-S-10B

Manual Integrations
 APPROVED

Sohil
 10/8/2018 12:21:14 PM

Quant Time: Oct 05 15:21:45 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 22 13:32:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	93660	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	382224	20.00	ng	-0.01
38) Acenaphthene-d10	9.73	164	184344	20.00	ng	-0.01
63) Phenanthrene-d10	11.21	188	328856	20.00	ng	0.00
75) Chrysene-d12	13.84	240	237077	20.00	ng	0.00
86) Perylene-d12	15.24	264	241916	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	472248	83.05	ng	0.01
7) Phenol-d6	6.36	99	633046	87.24	ng	0.00
23) Nitrobenzene-d5	7.26	82	412914	63.77	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	128792	70.87	ng	-0.02
44) 2-Fluorobiphenyl	9.06	172	788993	64.55	ng	-0.02
78) Terphenyl-d14	12.79	244	639758	66.23	ng	-0.01
Target Compounds						
10) Phenol	6.37	94	24737	3.010	ng	81
49) Dimethylphthalate	9.45	163	74846	5.582	ng	99
74) Fluoranthene	12.42	202	72981	4.159	ng	96
77) Pyrene	12.64	202	62631	3.686	ng	99
80) Benzo(a)anthracene	13.83	228	30425	2.131	ng	97
82) Chrysene	13.86	228	29803	2.106	ng	98
87) Benzo(b)fluoranthene	14.84	252	42585m	3.204	ng	
89) Benzo(a)pyrene	15.18	252	28032	2.340	ng	95

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

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