

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091118\
 Data File : BF108939.D
 Acq On : 11 Sep 2018 17:03
 Operator : JU/SJ
 Sample : J4838-07 2X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SUP-1A-C-090718-1

Manual Integrations
 APPROVED

Sohil
 9/12/2018 9:11:56 AM

Quant Time: Sep 11 17:51:58 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF090518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 05 18:03:31 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	181459	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	701884	20.00	ng	0.00
38) Acenaphthene-d10	9.77	164	312101	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	521169	20.00	ng	0.00
75) Chrysene-d12	13.87	240	418316	20.00	ng	0.00
86) Perylene-d12	15.28	264	495517	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	359381	32.81	ng	0.00
7) Phenol-d6	6.36	99	467157	33.15	ng	0.00
23) Nitrobenzene-d5	7.29	82	284494	25.37	ng	-0.01
41) 2,4,6-Tribromophenol	10.55	330	88551	29.72	ng	0.00
44) 2-Fluorobiphenyl	9.09	172	489273	26.74	ng	0.00
78) Terphenyl-d14	12.82	244	340879	19.00	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.48	163	160048	7.332	ng	99
70) Phenanthrene	11.27	178	52427	2.107	ng	98
74) Fluoranthene	12.45	202	84503	3.429	ng	98
77) Pyrene	12.68	202	75944	2.359	ng	98
87) Benzo(b)fluoranthene	14.88	252	60722m	2.254	ng	

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

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