

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090817\
 Data File : BF098373.D
 Acq On : 9 Sep 2017 4:09
 Operator : SJ/JU
 Sample : I5136-03
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BRICK-WALL

Manual Integrations
 APPROVED

Sohil
 9/11/2017 4:48:03 PM

Quant Time: Sep 09 05:38:21 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 05 17:28:20 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.68	152	101574	20.00	ng	-0.02
21) Naphthalene-d8	7.96	136	357876	20.00	ng	-0.02
38) Acenaphthene-d10	9.72	164	130900	20.00	ng	-0.02
63) Phenanthrene-d10	11.20	188	222996	20.00	ng	-0.02
75) Chrysene-d12	13.83	240	231990	20.00	ng	-0.03
86) Perylene-d12	15.24	264	160190	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	662421	99.20	ng	0.00
7) Phenol-d6	6.34	99	916001	100.96	ng	-0.01
23) Nitrobenzene-d5	7.25	82	515305	78.22	ng	-0.02
41) 2,4,6-Tribromophenol	10.51	330	113549	99.98	ng	-0.02
44) 2-Fluorobiphenyl	9.05	172	765965	85.97	ng	-0.02
78) Terphenyl-d14	12.79	244	604571	54.34	ng	-0.02
Target Compounds						
49) Dimethylphthalate	9.44	163	104832	10.956	ng	99
57) Fluorene	10.26	166	21867	2.416	ng	99
70) Phenanthrene	11.22	178	139119	11.708	ng	99
71) Anthracene	11.27	178	44116	3.660	ng	99
74) Fluoranthene	12.41	202	128417	10.354	ng	98
77) Pyrene	12.64	202	108985	5.744	ng	96
80) Benzo(a)anthracene	13.83	228	57396	4.128	ng	99
82) Chrysene	13.86	228	47618m	3.443	ng	
87) Benzo(b)fluoranthene	14.84	252	46096m	4.565	ng	
89) Benzo(a)pyrene	15.19	252	31010	3.469	ng	96

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

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