

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050117\
 Data File : BF094752.D
 Acq On : 1 May 2017 15:39
 Operator : SJ/MA
 Sample : I2929-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GF-TEST

Quant Time: May 02 02:42:30 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF041717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 17 14:59:23 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.76	152	118670	20.00	ng	0.00
21) Naphthalene-d8	7.25	136	450429	20.00	ng	0.00
38) Acenaphthene-d10	9.39	164	183547	20.00	ng	0.00
63) Phenanthrene-d10	11.21	188	303610	20.00	ng	0.00
75) Chrysene-d12	14.46	240	216814	20.00	ng	0.00
86) Perylene-d12	16.09	264	194498	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.32	112	685748	95.87	ng	0.02
7) Phenol-d6	5.40	99	830322	98.47	ng	0.00
23) Nitrobenzene-d5	6.41	82	501120	68.70	ng	0.00
41) 2,4,6-Tribromophenol	10.37	330	147482	102.73	ng	0.00
44) 2-Fluorobiphenyl	8.58	172	846329	67.84	ng	0.00
78) Terphenyl-d14	13.19	244	561617	69.24	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.08	163	312354	22.84	ng	99
70) Phenanthrene	11.24	178	71720	4.19	ng	96
74) Fluoranthene	12.70	202	151010	8.52	ng	97
77) Pyrene	12.98	202	133384	8.81	ng	98
80) Benzo(a)anthracene	14.45	228	61295	4.86	ng	99
82) Chrysene	14.49	228	52658	4.57	ng	96
85) Indeno(1,2,3-cd)pyrene	17.38	276	32311	2.95	ng	94
87) Benzo(b)fluoranthene	15.69	252	75214m	6.32	ng	
89) Benzo(a)pyrene	16.04	252	54432	4.92	ng	92
91) Benzo(g,h,i)perylene	17.76	276	32040	3.42	ng	96

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

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