

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090717\
 Data File : BF098330.D
 Acq On : 7 Sep 2017 22:58
 Operator : SJ/JU
 Sample : I5071-08
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 G58A-2.5-5

Manual Integrations
 APPROVED

Sohil
 9/8/2017 5:45:28 PM

Quant Time: Sep 08 03:00:16 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	6.69	152	100428	20.00	ng	-0.02
21) Naphthalene-d8	7.97	136	394823	20.00	ng	-0.01
38) Acenaphthene-d10	9.73	164	163844	20.00	ng	-0.01
63) Phenanthrene-d10	11.21	188	294529	20.00	ng	-0.01
75) Chrysene-d12	13.84	240	244245	20.00	ng	-0.02
86) Perylene-d12	15.26	264	167514	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	759532	115.04	ng	0.00
7) Phenol-d6	6.34	99	1023563	114.10	ng	-0.01
23) Nitrobenzene-d5	7.26	82	614080	84.49	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	170169	119.70	ng	-0.02
44) 2-Fluorobiphenyl	9.05	172	913729	81.93	ng	-0.02
78) Terphenyl-d14	12.80	244	721897	61.63	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.44	163	152445	12.729	ng	Qvalue 100
70) Phenanthrene	11.23	178	67248	4.285	ng	96
74) Fluoranthene	12.42	202	143883	8.783	ng	97
77) Pyrene	12.64	202	152544	7.636	ng	98
80) Benzo(a)anthracene	13.83	228	86755	5.926	ng	97
82) Chrysene	13.87	228	76481	5.252	ng	97
85) Indeno(1,2,3-cd)pyrene	16.61	276	22267	2.079	ng	93
87) Benzo(b)fluoranthene	14.86	252	88032m	8.337	ng	
88) Benzo(k)fluoranthene	14.88	252	20689m	2.086	ng	
89) Benzo(a)pyrene	15.20	252	44179	4.726	ng	96
91) Benzo(g,h,i)perylene	17.02	276	22448	2.827	ng	93

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

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