

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110317\
 Data File : BF100165.D
 Acq On : 4 Nov 2017 6:14
 Operator : SJ/JU
 Sample : I6265-01 2X
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-1

Manual Integrations
 APPROVED

Sohil
 11/4/2017 2:19:20 PM

Quant Time: Nov 04 06:58:28 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 03 15:46:10 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	95641	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	381344	20.00	ng	0.00
38) Acenaphthene-d10	9.80	164	163906	20.00	ng	0.00
63) Phenanthrene-d10	11.30	188	249406	20.00	ng	0.00
75) Chrysene-d12	13.95	240	158447	20.00	ng	0.00
86) Perylene-d12	15.42	264	156003	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	256879	42.17	ng	0.00
7) Phenol-d6	6.42	99	316277	43.79	ng	-0.01
23) Nitrobenzene-d5	7.34	82	179989	30.39	ng	-0.01
41) 2,4,6-Tribromophenol	10.60	330	58648	39.26	ng	0.00
44) 2-Fluorobiphenyl	9.12	172	353616	33.29	ng	0.00
78) Terphenyl-d14	12.87	244	201767	27.83	ng	0.00
Target Compounds						
10) Phenol	6.44	94	17089	2.155	ng	Qvalue 80
31) Naphthalene	8.07	128	53123	2.886	ng	99
49) Dimethylphthalate	9.53	163	35724	2.752	ng	# 98
70) Phenanthrene	11.32	178	55594	3.950	ng	99
74) Fluoranthene	12.52	202	72086	4.928	ng	96
77) Pyrene	12.75	202	65359	5.425	ng	98
80) Benzo(a)anthracene	13.94	228	31885m	3.174	ng	
82) Chrysene	13.98	228	31233	3.307	ng	97
85) Indeno(1,2,3-cd)pyrene	16.92	276	18732	2.161	ng	92
87) Benzo(b)fluoranthene	15.00	252	35986	3.653	ng	# 83
89) Benzo(a)pyrene	15.36	252	28450	3.215	ng	# 81
91) Benzo(g,h,i)perylene	17.37	276	17145	2.229	ng	# 84

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

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