

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120616\
 Data File : BF091465.D
 Acq On : 6 Dec 2016 18:16
 Operator : UM/SJ
 Sample : H5836-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP14-15-COMP2

Quant Time: Dec 07 00:51:30 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 07 00:12:14 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	123187	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	493934	20.00	ng	0.00
38) Acenaphthene-d10	9.88	164	272141	20.00	ng	0.00
63) Phenanthrene-d10	11.36	188	418636	20.00	ng	0.00
75) Chrysene-d12	13.99	240	338332	20.00	ng	0.00
86) Perylene-d12	15.43	264	302795	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	628822	83.39	ng	0.00
7) Phenol-d6	6.47	99	724327	78.03	ng	0.00
23) Nitrobenzene-d5	7.41	82	538264	61.07	ng	0.00
41) 2,4,6-Tribromophenol	10.66	330	227267	88.21	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	847970	56.42	ng	0.00
78) Terphenyl-d14	12.94	244	692572	44.49	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.48	94	50662	4.64	ng	98
49) Dimethylphthalate	9.59	163	134690	8.17	ng	# 96
70) Phenanthrene	11.38	178	82637	3.80	ng	96
74) Fluoranthene	12.56	202	167386	8.12	ng	98
77) Pyrene	12.79	202	215554	8.40	ng	96
80) Benzo(a)anthracene	13.98	228	163925	8.28	ng	96
82) Chrysene	14.01	228	113114m	5.78	ng	
85) Indeno(1,2,3-cd)pyrene	16.81	276	47956	2.67	ng	# 84
87) Benzo(b)fluoranthene	15.02	252	152549m	8.11	ng	
88) Benzo(k)fluoranthene	15.03	252	46199m	2.92	ng	
89) Benzo(a)pyrene	15.36	252	125383	7.42	ng	# 87
91) Benzo(g,h,i)perylene	17.23	276	50988	3.16	ng	95

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

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