

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120616\
 Data File : BF091468.D
 Acq On : 6 Dec 2016 19:41
 Operator : UM/SJ
 Sample : H5836-07
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP16-COMP3

Quant Time: Dec 07 01:02:27 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	126911	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	493191	20.00	ng	0.00
38) Acenaphthene-d10	9.88	164	258407	20.00	ng	0.00
63) Phenanthrene-d10	11.36	188	374359	20.00	ng	0.00
75) Chrysene-d12	13.99	240	276213	20.00	ng	0.00
86) Perylene-d12	15.43	264	244929	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	1033763	133.07	ng	0.01
7) Phenol-d6	6.48	99	1295010	135.42	ng	0.01
23) Nitrobenzene-d5	7.41	82	803560	91.31	ng	0.00
41) 2,4,6-Tribromophenol	10.66	330	269195	110.04	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	1059632	74.25	ng	0.00
78) Terphenyl-d14	12.95	244	794938	62.55	ng	0.01
Target Compounds						Qvalue
10) Phenol	6.49	94	76342	6.78	ng	86
49) Dimethylphthalate	9.60	163	196812	12.58	ng	# 99
70) Phenanthrene	11.38	178	126205	6.49	ng	99
71) Anthracene	11.43	178	39814	2.06	ng	96
74) Fluoranthene	12.57	202	263465	14.30	ng	99
77) Pyrene	12.79	202	233002m	11.12	ng	
80) Benzo(a)anthracene	13.98	228	148201	9.17	ng	95
82) Chrysene	14.01	228	111416m	6.97	ng	
85) Indeno(1,2,3-cd)pyrene	16.81	276	51320	3.51	ng	# 87
87) Benzo(b)fluoranthene	15.02	252	135782m	8.92	ng	
88) Benzo(k)fluoranthene	15.04	252	42346m	3.31	ng	
89) Benzo(a)pyrene	15.36	252	101584	7.43	ng	# 90
91) Benzo(g,h,i)perylene	17.23	276	48850	3.75	ng	# 92

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

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