

Data Path : Z:\HPCHEM1\BNA F\DATA\BF070517\
 Data File : BF096484.D
 Acq On : 5 Jul 2017 15:20
 Operator : SJ/MA
 Sample : I3871-03
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 POSTEX-51-20170622

Quant Time: Jul 05 17:04:02 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF070117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 05 13:27:10 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.71	152	118865	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	494887	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	199222	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	267469	20.00	ng	0.00
75) Chrysene-d12	13.86	240	224250	20.00	ng	0.00
86) Perylene-d12	15.27	264	187710	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	801591	99.54	ng	0.00
7) Phenol-d6	6.35	99	1058592	106.92	ng	0.00
23) Nitrobenzene-d5	7.27	82	554500	67.33	ng	0.00
41) 2,4,6-Tribromophenol	10.53	330	155191	99.73	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	959514	82.35	ng	0.00
78) Terphenyl-d14	12.82	244	541055	51.56	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.46	163	181057	12.17	ng	100
70) Phenanthrene	11.25	178	212514	14.70	ng	98
71) Anthracene	11.30	178	55064	3.74	ng	96
74) Fluoranthene	12.44	202	334670	23.61	ng	98
77) Pyrene	12.67	202	318727	17.71	ng	97
80) Benzo(a)anthracene	13.85	228	176004	13.55	ng	97
82) Chrysene	13.89	228	159820	11.75	ng	97
85) Indeno(1,2,3-cd)pyrene	16.63	276	66904	6.23	ng	99
87) Benzo(b)fluoranthene	14.87	252	168494m	14.33	ng	
88) Benzo(k)fluoranthene	14.89	252	63739m	6.06	ng	
89) Benzo(a)pyrene	15.21	252	124081	11.82	ng	# 95
91) Benzo(a,h,i)perylene	17.03	276	57469	6.71	ng	96

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

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