

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050117\
 Data File : BF094754.D
 Acq On : 1 May 2017 16:33
 Operator : SJ/MA
 Sample : I2927-03
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-03-003-B

Quant Time: May 02 02:58:35 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF041717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 17 14:59:23 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.76	152	105718	20.00	ng	0.00
21) Naphthalene-d8	7.25	136	402578	20.00	ng	0.00
38) Acenaphthene-d10	9.39	164	177420	20.00	ng	0.00
63) Phenanthrene-d10	11.21	188	268740	20.00	ng	0.00
75) Chrysene-d12	14.47	240	207634	20.00	ng	0.00
86) Perylene-d12	16.09	264	170907	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.33	112	716022	112.37	ng	0.03
7) Phenol-d6	5.40	99	880885	117.26	ng	0.00
23) Nitrobenzene-d5	6.41	82	550139	84.38	ng	0.00
41) 2,4,6-Tribromophenol	10.37	330	150213	108.24	ng	0.00
44) 2-Fluorobiphenyl	8.58	172	919678	76.27	ng	0.00
78) Terphenyl-d14	13.19	244	553767	71.29	ng	0.00
Target Compounds						Qvalue
10) Phenol	5.41	94	18642	2.02	ng	95
49) Dimethylphthalate	9.08	163	351570	26.59	ng	99
70) Phenanthrene	11.24	178	110612	7.31	ng	98
74) Fluoranthene	12.71	202	145207	9.26	ng	97
77) Pyrene	12.98	202	127358	8.78	ng	96
80) Benzo(a)anthracene	14.45	228	58456	4.84	ng	99
82) Chrysene	14.49	228	56049	5.08	ng	97
83) Bis(2-ethylhexyl)phthalate	14.51	149	29340	3.44	ng	98
85) Indeno(1,2,3-cd)pyrene	17.39	276	23561	2.25	ng	93
87) Benzo(b)fluoranthene	15.69	252	63817m	6.10	ng	
88) Benzo(k)fluoranthene	15.72	252	22896m	2.31	ng	
89) Benzo(a)pyrene	16.04	252	43730	4.50	ng	# 90
91) Benzo(a,h,i)perylene	17.77	276	23373	2.84	ng	# 94

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

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