

Data Path : Z:\HPCHEM1\BNA F\DATA\BF070517\
 Data File : BF096521.D
 Acq On : 6 Jul 2017 9:19
 Operator : SJ/MA
 Sample : I4029-05
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-P4-NSW

Quant Time: Jul 06 11:30:11 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	6.72	152	97453	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	354385	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	130037	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	231712	20.00	ng	0.00
75) Chrysene-d12	13.86	240	183771	20.00	ng	0.00
86) Perylene-d12	15.27	264	128139	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	777149	117.70	ng	0.00
7) Phenol-d6	6.35	99	831446	102.43	ng	0.00
23) Nitrobenzene-d5	7.27	82	431996	73.25	ng	0.00
41) 2,4,6-Tribromophenol	10.54	330	116966	115.15	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	681896	89.66	ng	0.00
78) Terphenyl-d14	12.82	244	503405	58.53	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.36	94	39579	4.28	ng	99
49) Dimethylphthalate	9.46	163	69678	7.18	ng	99
70) Phenanthrene	11.25	178	126642	10.11	ng	98
71) Anthracene	11.30	178	40194	3.15	ng	97
74) Fluoranthene	12.44	202	316297	25.76	ng	99
77) Pyrene	12.67	202	302053	20.48	ng	100
80) Benzo(a)anthracene	13.86	228	176342	16.57	ng	100
82) Chrysene	13.89	228	157683	14.15	ng	97
85) Indeno(1,2,3-cd)pyrene	16.64	276	74052	8.42	ng	95
87) Benzo(b)fluoranthene	14.88	252	182854m	22.77	ng	
88) Benzo(k)fluoranthene	14.90	252	58909m	8.20	ng	
89) Benzo(a)pyrene	15.22	252	127691	17.81	ng	95
90) Dibenzo(a,h)anthracene	16.64	278	18563	3.29	ng	# 84
91) Benzo(a,h,i)perylene	17.04	276	69317	11.86	ng	96

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

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