

Data Path : Z:\HPCHEM1\BNA F\DATA\BF070517\
 Data File : BF096501.D
 Acq On : 6 Jul 2017 00:07
 Operator : SJ/MA
 Sample : I3976-01
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B9-0-2

Quant Time: Jul 06 04:17:21 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	107901	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	417291	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	148458	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	207294	20.00	ng	-0.01
75) Chrysene-d12	13.86	240	221805	20.00	ng	-0.01
86) Perylene-d12	15.27	264	171473	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	633204	86.62	ng	0.00
7) Phenol-d6	6.35	99	770616	85.74	ng	0.00
23) Nitrobenzene-d5	7.27	82	441108	63.52	ng	-0.01
41) 2,4,6-Tribromophenol	10.53	330	79394	68.46	ng	-0.01
44) 2-Fluorobiphenyl	9.07	172	730286	84.11	ng	0.00
78) Terphenyl-d14	12.81	244	492023	47.40	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.46	163	102827	9.28	ng	98
70) Phenanthrene	11.25	178	81975	7.32	ng	98
74) Fluoranthene	12.43	202	138352	12.59	ng	98
77) Pyrene	12.66	202	136336	7.66	ng	99
80) Benzo(a)anthracene	13.84	228	67696	5.27	ng	97
82) Chrysene	13.88	228	66710	4.96	ng	97
85) Indeno(1,2,3-cd)pyrene	16.62	276	21506	2.03	ng	97
87) Benzo(b)fluoranthene	14.87	252	73464m	6.84	ng	
89) Benzo(a)pyrene	15.20	252	44433	4.63	ng	# 86
91) Benzo(g,h,i)perylene	17.03	276	25411	3.25	ng	97

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

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