

Data Path : Z:\HPCHEM1\BNA F\DATA\BF070517\
 Data File : BF096476.D
 Acq On : 5 Jul 2017 11:41
 Operator : SJ/MA
 Sample : I4031-01 5X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-02-063017

Quant Time: Jul 05 13:46:54 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	112349	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	457142	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	184003	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	265003	20.00	ng	0.00
75) Chrysene-d12	13.86	240	205872	20.00	ng	0.00
86) Perylene-d12	15.27	264	171686	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	207516	27.26	ng	0.00
7) Phenol-d6	6.34	99	218326	23.33	ng	-0.02
23) Nitrobenzene-d5	7.27	82	118523	15.58	ng	-0.01
41) 2,4,6-Tribromophenol	10.53	330	37219	25.90	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	234608	21.80	ng	-0.01
78) Terphenyl-d14	12.81	244	136852	14.20	ng	0.00
Target Compounds						Qvalue
48) Acenaphthylene	9.60	152	49117	2.64	ng	98
49) Dimethylphthalate	9.46	163	34563	2.52	ng	98
57) Fluorene	10.29	166	32274	3.02	ng	98
70) Phenanthrene	11.25	178	294873	20.58	ng	99
71) Anthracene	11.30	178	79697	5.46	ng	95
74) Fluoranthene	12.44	202	347663	24.75	ng	99
77) Pyrene	12.66	202	346448	20.97	ng	99
80) Benzo(a)anthracene	13.85	228	135444m	11.36	ng	
82) Chrysene	13.89	228	134400	10.77	ng	99
85) Indeno(1,2,3-cd)pyrene	16.63	276	46228	4.69	ng	96
87) Benzo(b)fluoranthene	14.87	252	134948m	12.54	ng	
88) Benzo(k)fluoranthene	14.89	252	48506m	5.04	ng	
89) Benzo(a)pyrene	15.21	252	100597	10.47	ng	97
91) Benzo(a,h,i)perylene	17.03	276	53268	6.80	ng	95

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

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