

Data Path : Z:\HPCHEM1\BNA F\DATA\BF070517\
 Data File : BF096502.D
 Acq On : 6 Jul 2017 00:35
 Operator : SJ/MA
 Sample : I3976-02
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B9-2-4

Quant Time: Jul 06 04:20:41 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	121420	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	472338	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	185962	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	255668	20.00	ng	0.00
75) Chrysene-d12	13.86	240	237421	20.00	ng	0.00
86) Perylene-d12	15.27	264	167651	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	425437	51.72	ng	0.00
7) Phenol-d6	6.35	99	808750	79.97	ng	0.00
23) Nitrobenzene-d5	7.27	82	494248	62.88	ng	-0.01
41) 2,4,6-Tribromophenol	10.53	330	35253	24.27	ng	-0.01
44) 2-Fluorobiphenyl	9.07	172	791539	72.78	ng	0.00
78) Terphenyl-d14	12.81	244	501868	45.17	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.46	163	84224	6.07	ng	99
70) Phenanthrene	11.25	178	232218	16.80	ng	96
71) Anthracene	11.30	178	48564	3.45	ng	98
74) Fluoranthene	12.43	202	336188	24.81	ng	98
77) Pyrene	12.66	202	309573	16.24	ng	99
80) Benzo(a)anthracene	13.85	228	142770	10.38	ng	100
82) Chrysene	13.89	228	136118	9.45	ng	96
85) Indeno(1,2,3-cd)pyrene	16.63	276	40011	3.52	ng	99
87) Benzo(b)fluoranthene	14.87	252	124002m	11.80	ng	
88) Benzo(k)fluoranthene	14.89	252	41182m	4.38	ng	
89) Benzo(a)pyrene	15.21	252	82024	8.75	ng	# 91
91) Benzo(a,h,i)perylene	17.04	276	39376	5.15	ng	93

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

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