

Data Path : Z:\HPCHEM1\BNA F\DATA\BF052217\
 Data File : BF095346.D
 Acq On : 23 May 2017 10:24
 Operator : SJ/MA
 Sample : I3231-01 5X
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RO-72-11975

Quant Time: May 23 12:10:35 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 18 17:07:53 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	70089	20.00	ng	0.00
21) Naphthalene-d8	9.78	136	315900	20.00	ng	0.00
38) Acenaphthene-d10	12.61	164	135903	20.00	ng	-0.01
63) Phenanthrene-d10	15.01	188	190315	20.00	ng	-0.01
75) Chrysene-d12	18.68	240	193230	20.00	ng	-0.01
86) Perylene-d12	20.36	264	141536	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	0.00	112	0	0.00	ng	
7) Phenol-d6	7.55	99	17146m	3.40	ng	0.24
23) Nitrobenzene-d5	8.80	82	7843m	1.57	ng	0.13
41) 2,4,6-Tribromophenol	0.00	330	0	0.00	ng	
44) 2-Fluorobiphenyl	11.57	172	30694	3.25	ng	0.01
78) Terphenyl-d14	17.33	244	19882	2.27	ng	-0.02
Target Compounds						Qvalue
31) Naphthalene	9.82	128	109279	6.88	ng	99
37) 2-Methylnaphthalene	10.97	142	37046	3.56	ng	96
51) Acenaphthene	12.67	154	73888	8.55	ng	99
54) Dibenzofuran	12.97	168	88436	7.39	ng	95
57) Fluorene	13.51	166	111646	10.73	ng	97
70) Phenanthrene	15.05	178	645335	59.02	ng	98
71) Anthracene	15.14	178	192490	17.23	ng	96
72) Carbazole	15.44	167	57838	5.69	ng	98
74) Fluoranthene	16.79	202	567380	50.58	ng	99
77) Pyrene	17.09	202	479455	33.18	ng	96
80) Benzo(a)anthracene	18.67	228	207491	18.45	ng	96
82) Chrysene	18.71	228	206889	19.03	ng	98
85) Indeno(1,2,3-cd)pyrene	21.71	276	56510	6.40	ng	# 90
87) Benzo(b)fluoranthene	19.95	252	172493m	19.72	ng	
88) Benzo(k)fluoranthene	19.98	252	72834m	8.63	ng	
89) Benzo(a)pyrene	20.31	252	126658	15.69	ng	97
91) Benzo(a,h,i)perylene	22.09	276	57800	8.84	ng	95

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

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