

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051215\
 Data File : BF079158.D
 Acq On : 12 May 2015 15:27
 Operator : TP/IZ
 Sample : G2155-07
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-44AS

Quant Time: May 13 11:40:56 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 13 00:53:14 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.28	152	65275	20.00	ng	0.00
21) Naphthalene-d8	8.86	136	283848	20.00	ng	0.00
38) Acenaphthene-d10	11.03	164	137345	20.00	ng	-0.01
63) Phenanthrene-d10	12.88	188	267492	20.00	ng	0.00
75) Chrysene-d12	16.22	240	273774	20.00	ng	0.02
86) Perylene-d12	18.10	264	300498	20.00	ng	0.08
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	386164	101.84	ng	0.01
7) Phenol-d6	6.84	99	486922	97.14	ng	0.00
23) Nitrobenzene-d5	7.98	82	294448	59.62	ng	0.00
41) 2,4,6-Tribromophenol	12.01	330	145772	98.11	ng	0.00
44) 2-Fluorobiphenyl	10.20	172	594324	60.29	ng	0.00
78) Terphenyl-d14	14.88	244	673881	58.93	ng	0.00
Target Compounds						Qvalue
32) Benzoic acid	8.51	122	6273	8.46	ng	94
49) Dimethylphthalate	10.71	163	35663	3.54	ng	100
70) Phenanthrene	12.90	178	273471	18.27	ng	100
71) Anthracene	12.97	178	66835	4.55	ng	98
72) Carbazole	13.18	167	31437	2.25	ng	99
74) Fluoranthene	14.39	202	403556	25.88	ng	95
77) Pyrene	14.67	202	362674	22.99	ng	98
80) Benzo(a)anthracene	16.21	228	196346	13.10	ng	97
82) Chrysene	16.25	228	172681	11.98	ng	98
85) Indeno(1,2,3-cd)pyrene	19.59	276	150898	8.21	ng	# 100
87) Benzo(b)fluoranthene	17.61	252	236478m	14.38	ng	
88) Benzo(k)fluoranthene	17.65	252	96780m	6.08	ng	
89) Benzo(a)pyrene	18.03	252	183126	12.09	ng	# 95
90) Dibenzo(a,h)anthracene	19.62	278	51791	3.22	ng	95
91) Benzo(a,h,i)perylene	20.02	276	142851	8.86	ng	95

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

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