

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051215\
 Data File : BF079194.D
 Acq On : 13 May 2015 11:21
 Operator : TP/IZ
 Sample : G2155-06DL 10X
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-48ASDL

Quant Time: May 13 13:46:43 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 13:12:37 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.28	152	51488	20.00	ng	0.00
21) Naphthalene-d8	8.86	136	213001	20.00	ng	0.00
38) Acenaphthene-d10	11.04	164	104886	20.00	ng	0.00
63) Phenanthrene-d10	12.88	188	199176	20.00	ng	0.00
75) Chrysene-d12	16.19	240	219717	20.00	ng	0.02
86) Perylene-d12	18.01	264	210746	20.00	ng	0.14
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	29739	9.94	ng	0.00
7) Phenol-d6	6.84	99	42888	10.85	ng	0.00
23) Nitrobenzene-d5	7.98	82	21551	5.81	ng	0.00
41) 2,4,6-Tribromophenol	12.02	330	9745	8.59	ng	0.00
44) 2-Fluorobiphenyl	10.20	172	45408	6.03	ng	0.00
78) Terphenyl-d14	14.88	244	50013	5.45	ng	0.00
Target Compounds						Qvalue
51) Acenaphthene	11.07	154	28319	4.45	ng	94
54) Dibenzofuran	11.29	168	25606	2.78	ng	96
57) Fluorene	11.71	166	22079	2.92	ng	96
70) Phenanthrene	12.91	178	464280	41.67	ng	99
71) Anthracene	12.97	178	92489	8.46	ng	99
72) Carbazole	13.18	167	23816	2.29	ng	98
74) Fluoranthene	14.39	202	564198	48.59	ng	99
77) Pyrene	14.67	202	536729	42.39	ng	99
80) Benzo(a)anthracene	16.18	228	225994	18.78	ng	99
82) Chrysene	16.23	228	219998m	19.01	ng	
85) Indeno(1,2,3-cd)pyrene	19.49	276	100616m	6.82	ng	
87) Benzo(b)fluoranthene	17.54	252	264634m	22.94	ng	
88) Benzo(k)fluoranthene	17.58	252	53213m	4.77	ng	
89) Benzo(a)pyrene	17.94	252	167130	15.74	ng	# 94
91) Benzo(a,h,i)perylene	19.92	276	96859	8.56	ng	97

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

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