

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050715\
 Data File : BF079011.D
 Acq On : 7 May 2015 22:29
 Operator : TP/IZ
 Sample : G2084-08
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 8

Manual Integrations
 APPROVED

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 5/8/2015 3:51:19 PM

Quant Time: May 08 11:40:22 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 08 00:46:40 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.35	152	65654	20.00	ng	0.01
21) Naphthalene-d8	8.92	136	274411	20.00	ng	0.01
38) Acenaphthene-d10	11.10	164	135145	20.00	ng	0.00
63) Phenanthrene-d10	12.95	188	275262	20.00	ng	0.01
75) Chrysene-d12	16.25	240	313970	20.00	ng	-0.02
86) Perylene-d12	18.03	264	321492	20.00	ng	-0.10
System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	338087	88.64	ng	0.01
7) Phenol-d6	6.90	99	472733	93.77	ng	0.01
23) Nitrobenzene-d5	8.03	82	266356	55.79	ng	0.00
41) 2,4,6-Tribromophenol	12.08	330	126378	86.44	ng	0.00
44) 2-Fluorobiphenyl	10.26	172	522013	53.81	ng	0.00
78) Terphenyl-d14	14.95	244	679335	51.80	ng	0.01
Target Compounds						Qvalue
49) Dimethylphthalate	10.77	163	27119	2.73	ng	99
70) Phenanthrene	12.97	178	78269	5.08	ng	99
74) Fluoranthene	14.46	202	182151	11.35	ng	96
77) Pyrene	14.74	202	153285	8.47	ng	97
80) Benzo(a)anthracene	16.24	228	115273m	6.70	ng	
82) Chrysene	16.29	228	143558m	8.68	ng	
85) Indeno(1,2,3-cd)pyrene	19.54	276	86002	4.08	ng	# 100
87) Benzo(b)fluoranthene	17.58	252	198878m	11.30	ng	
88) Benzo(k)fluoranthene	17.60	252	63519m	3.73	ng	
89) Benzo(a)pyrene	17.97	252	102613	6.33	ng	# 96
91) Benzo(g,h,i)perylene	19.99	276	79160	4.59	ng	98

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

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