

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
 Data File : BF078507.D
 Acq On : 14 Apr 2015 19:02
 Operator : TP/IZ
 Sample : G1828-03 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HS-SB06A

Quant Time: Apr 15 01:46:12 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 15 01:10:59 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	20008	20.00	ng	0.00
21) Naphthalene-d8	8.39	136	84701	20.00	ng	0.00
38) Acenaphthene-d10	10.54	164	46029	20.00	ng	0.00
63) Phenanthrene-d10	12.37	188	96873	20.00	ng	0.00
75) Chrysene-d12	15.62	240	113241	20.00	ng	-0.03
86) Perylene-d12	17.35	264	119041	20.00	ng	-0.14
System Monitoring Compounds						
5) 2-Fluorophenol	5.06	112	22206	18.30	ng	0.01
7) Phenol-d6	6.41	99	31453	20.68	ng	0.00
23) Nitrobenzene-d5	7.51	82	17395	12.45	ng	0.00
41) 2,4,6-Tribromophenol	11.52	330	8770	22.97	ng	0.00
44) 2-Fluorobiphenyl	9.72	172	39079	12.14	ng	-0.01
78) Terphenyl-d14	14.35	244	52667	10.51	ng	0.00
Target Compounds						Qvalue
48) Acenaphthylene	10.36	152	11788	2.46	ng	98
70) Phenanthrene	12.39	178	53668	9.58	ng	99
71) Anthracene	12.46	178	38050	6.89	ng	98
74) Fluoranthene	13.85	202	192725	32.61	ng	97
77) Pyrene	14.13	202	173579	24.42	ng	# 96
80) Benzo(a)anthracene	15.61	228	137435	20.40	ng	97
82) Chrysene	15.66	228	114066	18.02	ng	99
85) Indeno(1,2,3-cd)pyrene	18.58	276	61742m	8.60	ng	
87) Benzo(b)fluoranthene	16.90	252	176619	24.01	ng	# 95
88) Benzo(k)fluoranthene	16.94	252	37969m	5.81	ng	
89) Benzo(a)pyrene	17.28	252	91218	14.21	ng	# 94
91) Benzo(g,h,i)perylene	18.93	276	59158m	9.18	ng	

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

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