

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
 Data File : BF078511.D
 Acq On : 14 Apr 2015 20:57
 Operator : TP/IZ
 Sample : G1828-08
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HS-SB02B

Quant Time: Apr 15 02:04:30 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	19242	20.00	ng	0.00
21) Naphthalene-d8	8.39	136	83173	20.00	ng	0.00
38) Acenaphthene-d10	10.54	164	43946	20.00	ng	0.00
63) Phenanthrene-d10	12.36	188	85029	20.00	ng	0.00
75) Chrysene-d12	15.63	240	108278	20.00	ng	-0.02
86) Perylene-d12	17.37	264	118386	20.00	ng	-0.11
System Monitoring Compounds						
5) 2-Fluorophenol	5.06	112	26961	23.10	ng	0.01
7) Phenol-d6	6.41	99	55882	38.21	ng	0.00
23) Nitrobenzene-d5	7.51	82	37574	27.38	ng	0.00
41) 2,4,6-Tribromophenol	11.52	330	21592	59.24	ng	0.00
44) 2-Fluorobiphenyl	9.74	172	79072	25.72	ng	0.00
78) Terphenyl-d14	14.35	244	317717	66.30	ng	0.00
Target Compounds						
48) Acenaphthylene	10.36	152	12659	2.77	ng	# 90
70) Phenanthrene	12.40	178	31035	6.31	ng	# 89
71) Anthracene	12.46	178	13265	2.74	ng	# 76
74) Fluoranthene	13.86	202	103044	19.87	ng	94
77) Pyrene	14.14	202	123480	18.17	ng	# 96
80) Benzo(a)anthracene	15.62	228	124993	19.40	ng	96
82) Chrysene	15.66	228	105684m	17.46	ng	
85) Indeno(1,2,3-cd)pyrene	18.62	276	204393m	29.76	ng	
87) Benzo(b)fluoranthene	16.93	252	269642m	36.85	ng	
88) Benzo(k)fluoranthene	16.95	252	52068m	8.01	ng	
89) Benzo(a)pyrene	17.30	252	224354	35.14	ng	# 92
90) Dibenzo(a,h)anthracene	18.63	278	39966m	6.25	ng	
91) Benzo(g,h,i)perylene	18.97	276	217268m	33.90	ng	

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

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