

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011215\
 Data File : BF076830.D
 Acq On : 13 Jan 2015 5:19
 Operator : IZ / TP
 Sample : G1030-08 5X
 Misc : S
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP3-SS-2(4-6)

Quant Time: Jan 13 10:39:54 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF010915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 12 10:43:48 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	31415	20.00	ng	0.00
21) Naphthalene-d8	11.00	136	123666	20.00	ng	0.00
38) Acenaphthene-d10	15.12	164	65237	20.00	ng	-0.01
63) Phenanthrene-d10	17.85	188	104358	20.00	ng	0.00
75) Chrysene-d12	21.35	240	108248	20.00	ng	0.00
86) Perylene-d12	23.02	264	101117	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.17	112	38794	20.61	ng	-0.01
7) Phenol-d6	7.49	99	52383	22.10	ng	-0.01
23) Nitrobenzene-d5	9.37	82	32749	14.72	ng	-0.01
41) 2,4,6-Tribromophenol	16.77	330	10508	19.30	ng	-0.01
44) 2-Fluorobiphenyl	13.64	172	57881	13.06	ng	-0.01
78) Terphenyl-d14	20.07	244	51307	10.41	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	11.04	128	58860	9.28	ng	97
37) 2-Methylnaphthalene	12.70	142	61869	13.90	ng	96
45) 1,1'-Biphenyl	13.84	154	11312	2.26	ng	99
48) Acenaphthylene	14.78	152	81975	12.00	ng	99
51) Acenaphthene	15.20	154	63527	16.46	ng	95
54) Dibenzofuran	15.63	168	51212	9.07	ng	98
57) Fluorene	16.32	166	86665	18.38	ng	98
70) Phenanthrene	17.90	178	1072861	164.97	ng	97
71) Anthracene	17.97	178	308547	48.85	ng	99
72) Carbazole	18.24	167	117575	18.99	ng	97
74) Fluoranthene	19.55	202	1895813	284.62	ng	95
77) Pvrene	19.83	202	1670842	214.52	ng	97
80) Benzo(a)anthracene	21.34	228	972483	141.02	ng	96
82) Chrysene	21.39	228	732489m	117.40	ng	
85) Indeno(1,2,3-cd)pyrene	24.16	276	407619	66.02	ng	# 100
87) Benzo(b)fluoranthene	22.61	252	948121m	135.37	ng	
88) Benzo(k)fluoranthene	22.62	252	360839m	58.94	ng	
89) Benzo(a)pyrene	22.96	252	675619m	112.25	ng	
90) Dibenzo(a,h)anthracene	24.17	278	96234m	17.55	ng	
91) Benzo(a,h,i)perylene	24.47	276	377769	67.50	ng	99

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

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