

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011215\
 Data File : BF076834.D
 Acq On : 13 Jan 2015 7:37
 Operator : IZ / TP
 Sample : G1030-14
 Misc : S
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP5-SS-2(4-6)

Quant Time: Jan 13 11:43:18 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF010915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jan 10 06:22:22 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	35227	20.00	ng	0.00
21) Naphthalene-d8	11.00	136	146784	20.00	ng	0.00
38) Acenaphthene-d10	15.13	164	74502	20.00	ng	0.00
63) Phenanthrene-d10	17.87	188	118677	20.00	ng	0.01
75) Chrysene-d12	21.36	240	120381	20.00	ng	0.01
86) Perylene-d12	23.03	264	112933	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.17	112	190313	88.86	ng	-0.01
7) Phenol-d6	7.49	99	236637	87.70	ng	-0.01
23) Nitrobenzene-d5	9.37	82	147160	54.55	ng	-0.01
41) 2,4,6-Tribromophenol	16.77	330	53371	83.61	ng	-0.01
44) 2-Fluorobiphenyl	13.65	172	275697	53.85	ng	0.00
78) Terphenyl-d14	20.08	244	239922	42.93	ng	0.01
Target Compounds						Qvalue
31) Naphthalene	11.04	128	132035	17.42	ng	98
37) 2-Methylnaphthalene	12.70	142	37214	6.97	ng	96
45) 1,1'-Biphenyl	13.85	154	12886	2.21	ng	96
48) Acenaphthylene	14.78	152	90749	11.40	ng	97
49) Dimethylphthalate	14.70	163	30599	5.59	ng	# 93
51) Acenaphthene	15.20	154	120706	26.84	ng	100
54) Dibenzofuran	15.63	168	127558	19.60	ng	98
57) Fluorene	16.32	166	156335	28.46	ng	99
70) Phenanthrene	17.91	178	1973474	265.29	ng	98
71) Anthracene	17.98	178	506472	68.91	ng	99
72) Carbazole	18.25	167	246556	34.74	ng	99
74) Fluoranthene	19.55	202	2472281	322.26	ng	94
77) Pvrene	19.83	202	1818182	206.15	ng	# 89
80) Benzo(a)anthracene	21.35	228	1139210	145.80	ng	98
82) Chrysene	21.40	228	828564m	119.30	ng	
85) Indeno(1,2,3-cd)pyrene	24.20	276	457152	63.30	ng	# 100
87) Benzo(b)fluoranthene	22.62	252	1247312m	150.41	ng	
88) Benzo(k)fluoranthene	22.64	252	215476m	32.94	ng	
89) Benzo(a)pvrene	22.99	252	782941	113.94	ng	96
90) Dibenzo(a,h)anthracene	24.21	278	97865m	15.66	ng	
91) Benzo(q,h,i)perylene	24.51	276	414022	65.21	ng	99

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

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