

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011215\
 Data File : BF076835.D
 Acq On : 13 Jan 2015 8:11
 Operator : IZ / TP
 Sample : G1030-15
 Misc : S
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP5-SS-3(7-7.5)

Quant Time: Jan 13 11:53:50 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	34628	20.00	ng	0.00
21) Naphthalene-d8	11.00	136	139150	20.00	ng	0.00
38) Acenaphthene-d10	15.13	164	70819	20.00	ng	0.00
63) Phenanthrene-d10	17.87	188	116457	20.00	ng	0.01
75) Chrysene-d12	21.35	240	113910	20.00	ng	0.00
86) Perylene-d12	23.01	264	104533	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.17	112	220561	104.77	ng	-0.01
7) Phenol-d6	7.49	99	281163	106.01	ng	-0.01
23) Nitrobenzene-d5	9.37	82	170732	66.76	ng	-0.01
41) 2,4,6-Tribromophenol	16.77	330	61210	100.87	ng	-0.01
44) 2-Fluorobiphenyl	13.65	172	312617	64.24	ng	0.00
78) Terphenyl-d14	20.07	244	272134	51.46	ng	0.00

Target Compounds				Qvalue		
10) Phenol	7.51	94	6200	2.16	ng	84
31) Naphthalene	11.04	128	105023	14.62	ng	98
37) 2-Methylnaphthalene	12.70	142	26780	5.29	ng	94
48) Acenaphthylene	14.78	152	71301	9.42	ng	97
49) Dimethylphthalate	14.70	163	13454	2.59	ng	# 98
51) Acenaphthene	15.20	154	72934	17.06	ng	97
54) Dibenzofuran	15.63	168	72816	11.77	ng	100
57) Fluorene	16.32	166	101320	19.40	ng	96
70) Phenanthrene	17.90	178	1163664	159.41	ng	98
71) Anthracene	17.97	178	313280	43.44	ng	99
72) Carbazole	18.24	167	164455	23.62	ng	98
74) Fluoranthene	19.55	202	1607560	213.54	ng	97
77) Pyrene	19.82	202	1254691	150.34	ng	# 90
80) Benzo(a)anthracene	21.35	228	715868	96.82	ng	98
82) Chrysene	21.39	228	548811m	83.51	ng	
85) Indeno(1,2,3-cd)pyrene	24.15	276	309081	45.23	ng	# 100
87) Benzo(b)fluoranthene	22.61	252	775771m	101.07	ng	
88) Benzo(k)fluoranthene	22.63	252	205223m	33.89	ng	
89) Benzo(a)pyrene	22.95	252	523675	82.33	ng	# 92
90) Dibenz(a,h)anthracene	24.16	278	71752m	12.40	ng	
91) Benzo(g,h,i)perylene	24.46	276	276634	47.07	ng	95

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

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