

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011515\
 Data File : BF076886.D
 Acq On : 15 Jan 2015 16:14
 Operator : IZ / TP
 Sample : G1030-06 10X
 Misc : S
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP2-SS-3(7-8)

Quant Time: Jan 16 10:18:33 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 15 11:52:04 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	27716	20.00	ng	0.00
21) Naphthalene-d8	10.93	136	114352	20.00	ng	0.00
38) Acenaphthene-d10	15.07	164	60240	20.00	ng	-0.01
63) Phenanthrene-d10	17.82	188	94897	20.00	ng	0.00
75) Chrysene-d12	21.31	240	100278	20.00	ng	0.00
86) Perylene-d12	22.99	264	100902	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.10	112	22186	13.16	ng	0.01
7) Phenol-d6	7.42	99	26684	12.55	ng	0.00
23) Nitrobenzene-d5	9.32	82	17282	8.37	ng	0.00
41) 2,4,6-Tribromophenol	16.72	330	5334	10.91	ng	0.00
44) 2-Fluorobiphenyl	13.58	172	31802	8.03	ng	-0.01
78) Terphenyl-d14	20.04	244	30844	7.08	ng	0.00
Target Compounds						Qvalue
37) 2-Methylnaphthalene	12.63	142	20555	5.11	ng	97
48) Acenaphthylene	14.72	152	15073	2.43	ng	94
70) Phenanthrene	17.85	178	99512	16.82	ng	99
71) Anthracene	17.92	178	30960	5.36	ng	95
74) Fluoranthene	19.49	202	265987	43.86	ng	98
77) Pyrene	19.77	202	231745	33.72	ng	# 96
80) Benzo(a)anthracene	21.29	228	131642	20.92	ng	100
82) Chrysene	21.34	228	93722	16.33	ng	96
85) Indeno(1,2,3-cd)pyrene	24.12	276	53606	8.82	ng	# 86
87) Benzo(b)fluoranthene	22.56	252	127120m	18.71	ng	
88) Benzo(k)fluoranthene	22.57	252	50656m	8.53	ng	
89) Benzo(a)pyrene	22.92	252	95911	16.00	ng	# 95
90) Dibenzo(a,h)anthracene	24.14	278	12216m	2.22	ng	
91) Benzo(g,h,i)perylene	24.43	276	51670	9.45	ng	96

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

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