

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
 Data File : BF078519.D
 Acq On : 15 Apr 2015 00:46
 Operator : TP/IZ
 Sample : G1828-18
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HS-SB101

Quant Time: Apr 15 02:36:02 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	6.81	152	20153	20.00	ng	0.01
21) Naphthalene-d8	8.39	136	85738	20.00	ng	0.00
38) Acenaphthene-d10	10.54	164	43891	20.00	ng	0.00
63) Phenanthrene-d10	12.37	188	91520	20.00	ng	0.00
75) Chrysene-d12	15.62	240	106095	20.00	ng	-0.03
86) Perylene-d12	17.28	264	115145	20.00	ng	-0.21
System Monitoring Compounds						
5) 2-Fluorophenol	5.06	112	162442	132.90	ng	0.01
7) Phenol-d6	6.41	99	215356	140.59	ng	0.00
23) Nitrobenzene-d5	7.51	82	123944	87.62	ng	0.00
41) 2,4,6-Tribromophenol	11.52	330	56507	155.23	ng	0.00
44) 2-Fluorobiphenyl	9.74	172	249800	81.35	ng	0.00
78) Terphenyl-d14	14.35	244	310137	66.05	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.43	94	5256	2.86	ng	79
48) Acenaphthylene	10.36	152	12002	2.63	ng	96
49) Dimethylphthalate	10.25	163	12565	3.81	ng	97
70) Phenanthrene	12.39	178	51359	9.70	ng	99
71) Anthracene	12.46	178	35413	6.79	ng	99
74) Fluoranthene	13.85	202	175027	31.35	ng	99
77) Pyrene	14.13	202	156201	23.45	ng	# 93
80) Benzo(a)anthracene	15.61	228	111600	17.68	ng	97
82) Chrysene	15.66	228	73855	12.45	ng	95
85) Indeno(1,2,3-cd)pyrene	18.49	276	78421m	11.65	ng	
87) Benzo(b)fluoranthene	16.87	252	137830m	19.37	ng	
88) Benzo(k)fluoranthene	16.89	252	46729m	7.39	ng	
89) Benzo(a)pyrene	17.22	252	109640	17.65	ng	# 95
90) Dibenzo(a,h)anthracene	18.50	278	18299m	2.94	ng	
91) Benzo(g,h,i)perylene	18.83	276	74739m	11.99	ng	

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

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