

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
 Data File : BF078520.D
 Acq On : 15 Apr 2015 1:15
 Operator : TP/IZ
 Sample : G1828-06
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HS-SB03B

Quant Time: Apr 15 03:19:44 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	6.80	152	19176	20.00	ng	0.00
21) Naphthalene-d8	8.39	136	84201	20.00	ng	0.00
38) Acenaphthene-d10	10.54	164	45553	20.00	ng	0.00
63) Phenanthrene-d10	12.37	188	94371	20.00	ng	0.00
75) Chrysene-d12	15.62	240	114907	20.00	ng	-0.03
86) Perylene-d12	17.27	264	111844	20.00	ng	-0.22

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units
5) 2-Fluorophenol	0.00	112	0	0.00	ng
7) Phenol-d6	0.00	99	0	0.00	ng
23) Nitrobenzene-d5	0.00	82	0	0.00	ng
41) 2,4,6-Tribromophenol	0.00	330	0	0.00	ng
44) 2-Fluorobiphenyl	0.00	172	0	0.00	ng
78) Terphenyl-d14	0.00	244	0	0.00	ng

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
10) Phenol	6.44	94	3987m	2.28	ng	
49) Dimethylphthalate	10.25	163	14953	4.37	ng	# 94
70) Phenanthrene	12.39	178	21520	3.94	ng	99
74) Fluoranthene	13.85	202	64897	11.27	ng	96
77) Pyrene	14.13	202	58538	8.12	ng	# 92
80) Benzo(a)anthracene	15.61	228	27592	4.04	ng	98
82) Chrysene	15.65	228	28781	4.48	ng	97
85) Indeno(1,2,3-cd)pyrene	18.46	276	20406m	2.80	ng	
87) Benzo(b)fluoranthene	16.86	252	38046m	5.50	ng	
88) Benzo(k)fluoranthene	16.88	252	12859m	2.09	ng	
89) Benzo(a)pyrene	17.21	252	28068	4.65	ng	97
91) Benzo(g,h,i)perylene	18.80	276	19964m	3.30	ng	

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

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