

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
 Data File : BF078521.D
 Acq On : 15 Apr 2015 1:44
 Operator : TP/IZ
 Sample : G1823-11
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-6

Quant Time: Apr 15 03:23:23 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	20039	20.00	ng	0.00
21) Naphthalene-d8	8.39	136	91812	20.00	ng	0.00
38) Acenaphthene-d10	10.54	164	48973	20.00	ng	0.00
63) Phenanthrene-d10	12.37	188	102870	20.00	ng	0.00
75) Chrysene-d12	15.62	240	121827	20.00	ng	-0.03
86) Perylene-d12	17.30	264	129175	20.00	ng	-0.18

System Monitoring Compounds						
5) 2-Fluorophenol	5.05	112	124265	102.24	ng	0.00
7) Phenol-d6	6.41	99	168191	110.42	ng	0.00
23) Nitrobenzene-d5	7.51	82	95013	62.73	ng	0.00
41) 2,4,6-Tribromophenol	11.52	330	46626	114.80	ng	0.00
44) 2-Fluorobiphenyl	9.74	172	214590	62.63	ng	0.00
78) Terphenyl-d14	14.35	244	280909	52.10	ng	0.00

Target Compounds						Qvalue
10) Phenol	6.43	94	4141	2.27	ng	89
49) Dimethylphthalate	10.25	163	9830	2.67	ng	# 97
51) Acenaphthene	10.58	154	7078	2.47	ng	94
70) Phenanthrene	12.39	178	91096	15.31	ng	98
71) Anthracene	12.46	178	26143	4.46	ng	99
72) Carbazole	12.67	167	11605	2.11	ng	100
74) Fluoranthene	13.86	202	201126	32.05	ng	90
77) Pyrene	14.13	202	185406	24.24	ng	# 92
80) Benzo(a)anthracene	15.61	228	119557	16.49	ng	99
82) Chrysene	15.66	228	94914	13.94	ng	98
85) Indeno(1,2,3-cd)pyrene	18.50	276	60664m	7.85	ng	
87) Benzo(b)fluoranthene	16.87	252	145095	18.17	ng	# 95
88) Benzo(k)fluoranthene	16.90	252	42103m	5.93	ng	
89) Benzo(a)pyrene	17.23	252	89882	12.90	ng	# 95
90) Dibenzo(a,h)anthracene	18.53	278	15681m	2.25	ng	
91) Benzo(g,h,i)perylene	18.86	276	54792m	7.83	ng	

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

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