

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
 Data File : BF078522.D
 Acq On : 15 Apr 2015 2:13
 Operator : TP/IZ
 Sample : G1823-01 5X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-1

Quant Time: Apr 15 03:26:44 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	24624	20.00	ng	0.00
21) Naphthalene-d8	8.39	136	108945	20.00	ng	0.00
38) Acenaphthene-d10	10.54	164	57264	20.00	ng	0.00
63) Phenanthrene-d10	12.37	188	124374	20.00	ng	0.00
75) Chrysene-d12	15.62	240	142681	20.00	ng	-0.03
86) Perylene-d12	17.28	264	144268	20.00	ng	-0.21

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.06	112	23239	15.56	ng	0.01
7) Phenol-d6	6.41	99	31624	16.90	ng	0.00
23) Nitrobenzene-d5	7.51	82	17139	9.54	ng	0.00
41) 2,4,6-Tribromophenol	11.52	330	8131	17.12	ng	0.00
44) 2-Fluorobiphenyl	9.74	172	41919	10.46	ng	0.00
78) Terphenyl-d14	14.35	244	58349	9.24	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
48) Acenaphthylene	10.36	152	14460	2.43	ng	98
70) Phenanthrene	12.39	178	58513	8.13	ng	100
71) Anthracene	12.46	178	35532	5.01	ng	99
74) Fluoranthene	13.86	202	235120	30.99	ng	91
77) Pyrene	14.14	202	235277	26.27	ng	99
80) Benzo(a)anthracene	15.61	228	170039	20.03	ng	98
82) Chrysene	15.66	228	125836	15.78	ng	96
85) Indeno(1,2,3-cd)pyrene	18.49	276	75415m	8.33	ng	
87) Benzo(b)fluoranthene	16.87	252	194029m	21.76	ng	
88) Benzo(k)fluoranthene	16.89	252	59778m	7.54	ng	
89) Benzo(a)pyrene	17.22	252	120837	15.53	ng	97
90) Dibenzo(a,h)anthracene	18.50	278	16783m	2.15	ng	
91) Benzo(g,h,i)perylene	18.83	276	71147m	9.11	ng	

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

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