

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041615\
 Data File : BF078576.D
 Acq On : 16 Apr 2015 20:17
 Operator : TP/IZ
 Sample : G1855-06
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LS-SB15B(5.5-6)

Quant Time: Apr 17 02:23:55 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 17 00:52:22 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.68	152	21068	20.00	ng	0.00
21) Naphthalene-d8	8.27	136	91617	20.00	ng	0.00
38) Acenaphthene-d10	10.42	164	47890	20.00	ng	0.00
63) Phenanthrene-d10	12.24	188	97383	20.00	ng	0.00
75) Chrysene-d12	15.50	240	105824	20.00	ng	0.01
86) Perylene-d12	17.26	264	108705	20.00	ng	0.07

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	4.92	112	178020	139.32	ng	0.02
7) Phenol-d6	6.31	99	238075	148.67	ng	0.01
23) Nitrobenzene-d5	7.39	82	130447	86.30	ng	0.00
41) 2,4,6-Tribromophenol	11.40	330	61579	155.04	ng	0.01
44) 2-Fluorobiphenyl	9.62	172	272779	81.42	ng	0.00
78) Terphenyl-d14	14.24	244	317565	67.81	ng	0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
20) 3+4-Methylphenols	7.28	107	7033	4.49	ng	98
49) Dimethylphthalate	10.14	163	21820	6.06	ng	# 97
70) Phenanthrene	12.27	178	38087	6.76	ng	100
71) Anthracene	12.33	178	13401	2.41	ng	97
74) Fluoranthene	13.74	202	54182	9.12	ng	92
77) Pyrene	14.00	202	52851	7.96	ng	# 90
80) Benzo(a)anthracene	15.49	228	30812	4.89	ng	93
82) Chrysene	15.53	228	25166m	4.25	ng	
85) Indeno(1,2,3-cd)pyrene	18.48	276	21225m	3.16	ng	
87) Benzo(b)fluoranthene	16.80	252	39840m	5.93	ng	
89) Benzo(a)pyrene	17.19	252	29626	5.05	ng	# 80
91) Benzo(g,h,i)perylene	18.81	276	21094m	3.58	ng	

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

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