

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042015\
 Data File : BF078664.D
 Acq On : 21 Apr 2015 3:19
 Operator : TP/IZ
 Sample : G1855-02
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LS-SB12B

Manual Integrations
 APPROVED

apatel
 4/21/2015 5:03:52 PM

Quant Time: Apr 21 04:01:21 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 21 03:43:19 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.86	152	37455	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	154264	20.00	ng	0.00
38) Acenaphthene-d10	11.20	164	78382	20.00	ng	-0.01
63) Phenanthrene-d10	13.93	188	153966	20.00	ng	-0.01
75) Chrysene-d12	17.79	240	159019	20.00	ng	-0.01
86) Perylene-d12	19.46	264	153928	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	3.87	112	210189	92.53	ng	0.00
7) Phenol-d6	5.36	99	278312	97.76	ng	0.00
23) Nitrobenzene-d5	6.80	82	177461	69.73	ng	0.00
41) 2,4,6-Tribromophenol	12.67	330	61417	94.48	ng	-0.01
44) 2-Fluorobiphenyl	10.03	172	323357	58.97	ng	0.00
78) Terphenyl-d14	16.46	244	253097	35.97	ng	-0.01
Target Compounds						
49) Dimethylphthalate	10.83	163	37539	6.37	ng	# 97
74) Fluoranthene	15.86	202	53816	5.73	ng	93
77) Pyrene	16.17	202	64042	6.42	ng	95
80) Benzo(a)anthracene	17.78	228	47892	5.06	ng	96
82) Chrysene	17.83	228	41080	4.62	ng	99
85) Indeno(1,2,3-cd)pyrene	20.61	276	47059	4.67	ng	# 89
87) Benzo(b)fluoranthene	19.06	252	75956m	7.98	ng	
88) Benzo(k)fluoranthene	19.09	252	28638m	3.39	ng	
89) Benzo(a)pyrene	19.41	252	55928	6.74	ng	# 95
91) Benzo(g,h,i)perylene	20.93	276	43142	5.18	ng	97

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

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