

Data Path : V:\HPCHEM1\BNA F\DATA\BF121515\
 Data File : BF083825.D
 Acq On : 15 Dec 2015 18:23
 Operator : UM/IZ
 Sample : G4739-15
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 K201-0-2

Quant Time: Dec 16 06:56:36 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 14 01:30:12 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	96334	20.00	ng	-0.02
21) Naphthalene-d8	8.19	136	379599	20.00	ng	-0.02
38) Acenaphthene-d10	9.94	164	145928	20.00	ng	-0.02
63) Phenanthrene-d10	11.42	188	235159	20.00	ng	-0.02
75) Chrysene-d12	14.07	240	196669	20.00	ng	-0.01
86) Perylene-d12	15.51	264	147924	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.52	112	682953	114.37	ng	-0.01
7) Phenol-d6	6.54	99	987625	130.60	ng	0.00
23) Nitrobenzene-d5	7.47	82	594395	87.60	ng	-0.01
41) 2,4,6-Tribromophenol	10.73	330	164200	110.43	ng	-0.02
44) 2-Fluorobiphenyl	9.26	172	823291	81.67	ng	-0.02
78) Terphenyl-d14	13.01	244	627958	68.68	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
6) Aniline	6.54	93	74914	7.29	ng	# 82
10) Phenol	6.56	94	22843	2.62	ng	# 63
31) Naphthalene	8.21	128	105294	5.32	ng	99
37) 2-Methylnaphthalene	8.90	142	61548	4.95	ng	97
48) Acenaphthylene	9.80	152	131982	8.22	ng	99
49) Dimethylphthalate	9.65	163	182944	16.08	ng	99
57) Fluorene	10.49	166	22081m	2.18	ng	
70) Phenanthrene	11.45	178	275133	22.02	ng	99
71) Anthracene	11.51	178	102077	7.87	ng	98
74) Fluoranthene	12.64	202	476847	36.94	ng	96
77) Pyrene	12.87	202	539556	34.71	ng	99
80) Benzo(a)anthracene	14.05	228	281774	25.20	ng	# 89
82) Chrysene	14.09	228	287075m	26.32	ng	
83) Bis(2-ethylhexyl)phthalate	14.05	149	53733	7.08	ng	# 98
85) Indeno(1,2,3-cd)pyrene	16.95	276	128753	12.02	ng	# 100
87) Benzo(b)fluoranthene	15.09	252	288283m	30.55	ng	
88) Benzo(k)fluoranthene	15.11	252	100216m	11.72	ng	
89) Benzo(a)pyrene	15.45	252	205880	24.45	ng	98
90) Dibenzo(a,h)anthracene	16.96	278	31161	3.72	ng	# 80
91) Benzo(a,h,i)perylene	17.38	276	138892	16.14	ng	99

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

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