

Data Path : V:\HPCHEM1\BNA F\DATA\BF121515\
 Data File : BF083822.D
 Acq On : 15 Dec 2015 16:59
 Operator : UM/IZ
 Sample : G4739-09
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 K206-18-19

Quant Time: Dec 16 04:44:27 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	94618	20.00	ng	-0.02
21) Naphthalene-d8	8.19	136	384226	20.00	ng	-0.02
38) Acenaphthene-d10	9.94	164	167210	20.00	ng	-0.02
63) Phenanthrene-d10	11.42	188	270829	20.00	ng	-0.02
75) Chrysene-d12	14.07	240	211967	20.00	ng	-0.01
86) Perylene-d12	15.51	264	167448	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	684390	116.69	ng	-0.01
7) Phenol-d6	6.54	99	972436	130.93	ng	0.00
23) Nitrobenzene-d5	7.46	82	560346	81.59	ng	-0.02
41) 2,4,6-Tribromophenol	10.73	330	201929	118.51	ng	-0.02
44) 2-Fluorobiphenyl	9.26	172	903816	76.96	ng	-0.02
78) Terphenyl-d14	13.01	244	748008	75.91	ng	-0.01
Target Compounds						Qvalue
31) Naphthalene	8.21	128	70267	3.51	ng	98
48) Acenaphthylene	9.80	152	218896	11.89	ng	99
49) Dimethylphthalate	9.65	163	156130	11.98	ng	99
51) Acenaphthene	9.97	154	73081	6.57	ng	98
54) Dibenzofuran	10.14	168	59371	4.03	ng	# 87
57) Fluorene	10.49	166	140721	12.12	ng	99
70) Phenanthrene	11.46	178	1041330	72.38	ng	99
71) Anthracene	11.51	178	481308	32.21	ng	100
72) Carbazole	11.65	167	56332	4.05	ng	# 92
74) Fluoranthene	12.65	202	1459981	98.21	ng	98
77) Pyrene	12.88	202	1532409	91.47	ng	97
80) Benzo(a)anthracene	14.05	228	650111m	53.94	ng	
82) Chrysene	14.09	228	553285m	47.06	ng	
85) Indeno(1,2,3-cd)pyrene	16.93	276	178657	15.47	ng	# 100
87) Benzo(b)fluoranthene	15.09	252	497246m	46.55	ng	
88) Benzo(k)fluoranthene	15.12	252	149915m	15.49	ng	
89) Benzo(a)pyrene	15.45	252	410847	43.10	ng	98
90) Dibenzo(a,h)anthracene	16.93	278	41449	4.37	ng	# 79
91) Benzo(a,h,i)perylene	17.37	276	179792	18.45	ng	99

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

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