

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091415\
 Data File : BF081608.D
 Acq On : 14 Sep 2015 15:10
 Operator : UM/IZ
 Sample : G3597-02 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GP-2(0-5)

Quant Time: Sep 15 01:49:58 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 14 13:51:56 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.23	152	43986	20.00	ng	0.00
21) Naphthalene-d8	11.12	136	176889	20.00	ng	-0.01
38) Acenaphthene-d10	15.26	164	85101	20.00	ng	-0.01
63) Phenanthrene-d10	18.77	188	167772	20.00	ng	0.00
75) Chrysene-d12	23.40	240	107816	20.00	ng	0.00
86) Perylene-d12	24.84	264	96721	20.00	ng	0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	63645	22.45	ng	0.00
7) Phenol-d6	7.63	99	93692	24.97	ng	-0.02
23) Nitrobenzene-d5	9.51	82	60717	16.56	ng	-0.03
41) 2,4,6-Tribromophenol	17.16	330	18313	24.17	ng	-0.02
44) 2-Fluorobiphenyl	13.76	172	110107	16.58	ng	-0.02
78) Terphenyl-d14	22.12	244	88686	19.15	ng	0.00

Target Compounds				Qvalue		
48) Acenaphthylene	14.92	152	88154	8.79	ng	98
49) Dimethylphthalate	14.82	163	25688	3.89	ng	# 96
57) Fluorene	16.57	166	13859	2.19	ng	91
70) Phenanthrene	18.82	178	367285	39.38	ng	98
71) Anthracene	18.94	178	84583	8.83	ng	99
72) Carbazole	19.42	167	23210	2.58	ng	93
74) Fluoranthene	21.45	202	773380	76.59	ng	89
77) Pyrene	21.59	202	34424	4.31	ng	95
80) Benzo(a)anthracene	23.38	228	394017	57.39	ng	95
82) Chrysene	23.43	228	346397	56.28	ng	94
85) Indeno(1,2,3-cd)pyrene	25.89	276	161473	35.76	ng	# 79
87) Benzo(b)fluoranthene	24.49	252	368614m	53.38	ng	
88) Benzo(k)fluoranthene	24.50	252	140524m	24.53	ng	
89) Benzo(a)pyrene	24.79	252	336231	57.46	ng	# 81
90) Dibenzo(a,h)anthracene	25.89	278	48305m	10.68	ng	
91) Benzo(g,h,i)perylene	26.20	276	148468	34.40	ng	# 72

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

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