

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091715\
 Data File : BF081685.D
 Acq On : 17 Sep 2015 18:53
 Operator : UM/IZ
 Sample : G3620-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-14(0-0.16)

Quant Time: Sep 18 07:16:23 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 15 14:03:59 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.17	152	44465	20.00	ng	-0.04
21) Naphthalene-d8	11.06	136	176060	20.00	ng	-0.06
38) Acenaphthene-d10	15.19	164	82081	20.00	ng	-0.07
63) Phenanthrene-d10	18.70	188	165495	20.00	ng	-0.06
75) Chrysene-d12	23.34	240	115745	20.00	ng	-0.04
86) Perylene-d12	24.78	264	74922	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	237737	82.95	ng	-0.02
7) Phenol-d6	7.59	99	331912	87.49	ng	-0.06
23) Nitrobenzene-d5	9.46	82	223640	61.29	ng	-0.07
41) 2,4,6-Tribromophenol	17.10	330	64068	87.66	ng	-0.07
44) 2-Fluorobiphenyl	13.71	172	384955	60.10	ng	-0.07
78) Terphenyl-d14	22.07	244	327921	65.95	ng	-0.04
Target Compounds						Qvalue
49) Dimethylphthalate	14.77	163	75637	11.86	ng	# 96
74) Fluoranthene	21.39	202	24577	2.47	ng	87
77) Pyrene	21.74	202	23734	2.77	ng	97
80) Benzo(a)anthracene	23.34	228	16039	2.18	ng	97
82) Chrysene	23.37	228	17821m	2.70	ng	
85) Indeno(1,2,3-cd)pyrene	25.82	276	10483	2.16	ng	# 78
87) Benzo(b)fluoranthene	24.45	252	23549m	4.40	ng	
89) Benzo(a)pyrene	24.73	252	13812m	3.05	ng	
91) Benzo(g,h,i)perylene	26.12	276	10053	3.01	ng	# 73

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

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