

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082515\
 Data File : BF081213.D
 Acq On : 25 Aug 2015 22:41
 Operator : UM/IZ
 Sample : G3407-10 5X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P004-BF003-01

Quant Time: Aug 26 01:31:44 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 24 13:58:19 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.53	152	69770	20.00	ng	-0.01
21) Naphthalene-d8	7.82	136	267004	20.00	ng	-0.01
38) Acenaphthene-d10	9.56	164	109576	20.00	ng	-0.02
63) Phenanthrene-d10	11.03	188	198408	20.00	ng	-0.01
75) Chrysene-d12	13.65	240	139488	20.00	ng	-0.01
86) Perylene-d12	14.99	264	113044	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.10	112	141311	30.46	ng	0.00
7) Phenol-d6	6.18	99	188066	31.07	ng	-0.01
23) Nitrobenzene-d5	7.10	82	114091	19.52	ng	-0.02
41) 2,4,6-Tribromophenol	10.34	330	26506	27.91	ng	-0.01
44) 2-Fluorobiphenyl	8.89	172	190280	22.75	ng	-0.01
78) Terphenyl-d14	12.61	244	119035	18.29	ng	-0.02
Target Compounds						Qvalue
49) Dimethylphthalate	9.29	163	40731	4.51	ng	100
70) Phenanthrene	11.05	178	46106	3.92	ng	100
74) Fluoranthene	12.23	202	59282	4.67	ng	97
77) Pyrene	12.45	202	49089m	4.33	ng	
80) Benzo(a)anthracene	13.64	228	22390	2.39	ng	93
82) Chrysene	13.67	228	24271m	2.88	ng	
87) Benzo(b)fluoranthene	14.63	252	28105m	3.51	ng	
89) Benzo(a)pyrene	14.93	252	19754	2.84	ng	# 87
91) Benzo(g,h,i)perylene	16.51	276	11602	2.11	ng	# 87

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

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