

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102115\
 Data File : BF082428.D
 Acq On : 22 Oct 2015 3:39
 Operator : UM/IZ
 Sample : G4116-07
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BRIDGE-14-01(0-2)

Quant Time: Oct 22 16:15:05 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 21 13:20:24 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	90360	20.00	ng	-0.01
21) Naphthalene-d8	8.06	136	337848	20.00	ng	-0.01
38) Acenaphthene-d10	9.82	164	140591	20.00	ng	0.00
63) Phenanthrene-d10	11.30	188	227982	20.00	ng	0.00
75) Chrysene-d12	13.96	240	210832	20.00	ng	-0.06
86) Perylene-d12	15.41	264	208540	20.00	ng	-0.13
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	733997	163.94	ng	0.01
7) Phenol-d6	6.42	99	939560	161.26	ng	0.00
23) Nitrobenzene-d5	7.34	82	533422	106.03	ng	-0.01
41) 2,4,6-Tribromophenol	10.61	330	161890	122.79	ng	-0.01
44) 2-Fluorobiphenyl	9.14	172	1009721	119.10	ng	0.00
78) Terphenyl-d14	12.89	244	716176	90.02	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.53	163	212794	21.49	ng	100
70) Phenanthrene	11.33	178	46540	4.16	ng	97
74) Fluoranthene	12.52	202	75454	6.29	ng	98
77) Pyrene	12.74	202	69603	5.56	ng	98
79) Butylbenzylphthalate	13.36	149	55416	9.71	ng	# 85
80) Benzo(a)anthracene	13.94	228	38681m	3.53	ng	
82) Chrysene	13.98	228	37007m	3.20	ng	
85) Indeno(1,2,3-cd)pyrene	16.80	276	21846	2.38	ng	94
87) Benzo(b)fluoranthene	15.01	252	50666m	3.99	ng	
88) Benzo(k)fluoranthene	15.02	252	21769m	2.04	ng	
89) Benzo(a)pyrene	15.35	252	35486	3.25	ng	# 88
91) Benzo(a,h,i)perylene	17.22	276	21797	2.51	ng	# 87

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

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