

Data Path : U:\HPCHEM1\BNA F\DATA\BF101915\
 Data File : BF082387.D
 Acq On : 20 Oct 2015 7:25
 Operator : UM/IZ
 Sample : G4046-01DL2 250X
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GOODLUCK902DL2

Manual Integrations
 APPROVED

MMdadoda
 10/20/2015 6:26:08 PM

Quant Time: Oct 20 08:50:46 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 12 01:46:21 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	83299	20.00	ng	-0.03
21) Naphthalene-d8	8.09	136	346134	20.00	ng	-0.03
38) Acenaphthene-d10	9.85	164	165438	20.00	ng	-0.03
63) Phenanthrene-d10	11.34	188	283576	20.00	ng	-0.03
75) Chrysene-d12	14.01	240	223607	20.00	ng	0.00
86) Perylene-d12	15.51	264	213886	20.00	ng	0.07
System Monitoring Compounds						
5) 2-Fluorophenol	0.00	112	0	0.00	ng	
7) Phenol-d6	6.48	99	441	0.08	ng	0.00
23) Nitrobenzene-d5	0.00	82	0	0.00	ng	
41) 2,4,6-Tribromophenol	10.64	330	209	0.13	ng	-0.03
44) 2-Fluorobiphenyl	9.17	172	3324	0.33	ng	-0.04
78) Terphenyl-d14	12.93	244	3151	0.37	ng	-0.03
Target Compounds						
57) Fluorene	10.40	166	29073	2.78	ng	97
70) Phenanthrene	11.36	178	544328	39.16	ng	99
71) Anthracene	11.42	178	241879	16.89	ng	99
72) Carbazole	11.58	167	82678	6.33	ng	97
74) Fluoranthene	12.56	202	612463	41.03	ng	98
77) Pyrene	12.79	202	318173	23.98	ng	97
80) Benzo(a)anthracene	14.00	228	76511	6.58	ng	99
82) Chrysene	14.04	228	80225m	6.54	ng	
87) Benzo(b)fluoranthene	15.09	252	39036m	3.00	ng	

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

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