

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100915\
 Data File : BF082150.D
 Acq On : 9 Oct 2015 17:39
 Operator : UM/IZ
 Sample : G3906-08
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 IM-MLA-(1.5-02)-R

Manual Integrations
 APPROVED

MMDadoda
 10/12/2015 2:43:47 PM

Quant Time: Oct 10 02:28:33 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 09 22:47:26 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	112531	20.00	ng	-0.01
21) Naphthalene-d8	8.14	136	449195	20.00	ng	0.00
38) Acenaphthene-d10	9.90	164	199165	20.00	ng	0.00
63) Phenanthrene-d10	11.38	188	342258	20.00	ng	0.00
75) Chrysene-d12	14.02	240	259391	20.00	ng	-0.01
86) Perylene-d12	15.46	264	217645	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	642044	103.57	ng	0.01
7) Phenol-d6	6.49	99	886442	113.64	ng	0.00
23) Nitrobenzene-d5	7.42	82	512941	76.12	ng	-0.01
41) 2,4,6-Tribromophenol	10.69	330	199901	99.10	ng	-0.01
44) 2-Fluorobiphenyl	9.22	172	970459	80.74	ng	0.00
78) Terphenyl-d14	12.97	244	746166	96.47	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.61	163	239164	17.40	ng	Qvalue # 97
70) Phenanthrene	11.41	178	67514	2.53	ng	99
74) Fluoranthene	12.59	202	86028	4.69	ng	88
77) Pyrene	12.82	202	77607	5.01	ng	98
80) Benzo(a)anthracene	14.01	228	32901m	2.36	ng	
82) Chrysene	14.05	228	31465m	2.59	ng	
87) Benzo(b)fluoranthene	15.04	252	32500m	2.62	ng	
89) Benzo(a)pyrene	15.40	252	24735	2.29	ng	# 91

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

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