

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101015\
 Data File : BF082212.D
 Acq On : 11 Oct 2015 9:09
 Operator : UM/IZ
 Sample : G3933-01
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-COMP-04

Manual Integrations
 APPROVED

MMDadoda
 10/12/2015 2:39:28 PM

Quant Time: Oct 12 04:41:02 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 02:59:50 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	96932	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	373775	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	156863	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	258077	20.00	ng	0.00
75) Chrysene-d12	14.01	240	253321	20.00	ng	0.00
86) Perylene-d12	15.45	264	204324	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	525998	109.52	ng	0.01
7) Phenol-d6	6.48	99	694454	111.11	ng	0.00
23) Nitrobenzene-d5	7.41	82	374203	67.23	ng	-0.01
41) 2,4,6-Tribromophenol	10.67	330	137393	93.40	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	702854	74.31	ng	-0.01
78) Terphenyl-d14	12.95	244	557750	58.34	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.60	163	143087	12.95	ng	99
70) Phenanthrene	11.39	178	28537	2.26	ng	98
74) Fluoranthene	12.58	202	65612	4.83	ng	97
77) Pyrene	12.81	202	79498	5.29	ng	99
80) Benzo(a)anthracene	14.00	228	48404m	3.67	ng	
82) Chrysene	14.04	228	44919m	3.23	ng	
85) Indeno(1,2,3-cd)pyrene	16.86	276	26909	2.44	ng	96
87) Benzo(b)fluoranthene	15.03	252	51186m	4.12	ng	
88) Benzo(k)fluoranthene	15.04	252	29493m	2.82	ng	
89) Benzo(a)pyrene	15.38	252	45691	4.28	ng	# 79
91) Benzo(g,h,i)perylene	17.29	276	28975	3.41	ng	96

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

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