

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101015\
 Data File : BF082205.D
 Acq On : 11 Oct 2015 5:40
 Operator : UM/IZ
 Sample : G3979-08
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 A16COMP

Manual Integrations
 APPROVED

MMDadoda
 10/12/2015 2:38:27 PM

Quant Time: Oct 12 04:09:14 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	85503	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	322176	20.00	ng	0.00
38) Acenaphthene-d10	9.87	164	140649	20.00	ng	-0.01
63) Phenanthrene-d10	11.36	188	230038	20.00	ng	-0.01
75) Chrysene-d12	14.00	240	191546	20.00	ng	-0.01
86) Perylene-d12	15.44	264	194403	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	531002	125.34	ng	0.01
7) Phenol-d6	6.48	99	675114	122.45	ng	0.00
23) Nitrobenzene-d5	7.41	82	387270	80.72	ng	-0.01
41) 2,4,6-Tribromophenol	10.67	330	145713	110.47	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	671014	79.12	ng	-0.01
78) Terphenyl-d14	12.95	244	472712	65.40	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.60	163	239890	24.22	ng	99
70) Phenanthrene	11.39	178	146312	12.98	ng	98
71) Anthracene	11.44	178	24060	2.07	ng	97
74) Fluoranthene	12.58	202	286250	23.64	ng	96
77) Pyrene	12.81	202	238804	21.01	ng	97
80) Benzo(a)anthracene	14.00	228	116147	11.65	ng	96
82) Chrysene	14.04	228	121393	11.56	ng	99
85) Indeno(1,2,3-cd)pyrene	16.85	276	81344	9.75	ng	96
87) Benzo(b)fluoranthene	15.03	252	183497m	15.51	ng	
88) Benzo(k)fluoranthene	15.04	252	82222m	8.27	ng	
89) Benzo(a)pyrene	15.38	252	118687	11.68	ng	98
91) Benzo(g,h,i)perylene	17.27	276	75705	9.36	ng	97

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

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