

Data Path : Z:\HPCHEM1\BNA F\DATA\BF090315\
 Data File : BF081412.D
 Acq On : 4 Sep 2015 9:57
 Operator : UM/IZ
 Sample : G3505-04
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P005-BF001-02

Manual Integrations
 APPROVED

MMDadoda
 9/4/2015 2:52:00 PM

Quant Time: Sep 04 13:26:54 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 01 17:35:35 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.36	152	62602	20.00	ng	-0.02
21) Naphthalene-d8	7.66	136	239267	20.00	ng	-0.02
38) Acenaphthene-d10	9.39	164	100885	20.00	ng	-0.02
63) Phenanthrene-d10	10.86	188	198032	20.00	ng	-0.02
75) Chrysene-d12	13.46	240	122674	20.00	ng	-0.03
86) Perylene-d12	14.78	264	103492	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	4.92	112	475703	117.90	ng	0.00
7) Phenol-d6	6.04	99	706769	132.33	ng	-0.01
23) Nitrobenzene-d5	6.95	82	447594	90.25	ng	-0.02
41) 2,4,6-Tribromophenol	10.18	330	105568	117.52	ng	-0.02
44) 2-Fluorobiphenyl	8.74	172	654152	83.09	ng	-0.02
78) Terphenyl-d14	12.44	244	488580	92.71	ng	-0.02
Target Compounds						
49) Dimethylphthalate	9.14	163	128075	16.34	ng	# 97
74) Fluoranthene	12.06	202	49243	4.13	ng	89
77) Pyrene	12.27	202	43423m	4.78	ng	
80) Benzo(a)anthracene	13.46	228	18414	2.36	ng	94
82) Chrysene	13.50	228	18780m	2.68	ng	
85) Indeno(1,2,3-cd)pyrene	15.85	276	10515	2.05	ng	# 89
87) Benzo(b)fluoranthene	14.45	252	27263m	3.69	ng	
89) Benzo(a)pyrene	14.73	252	16277m	2.60	ng	
91) Benzo(g,h,i)perylene	16.17	276	10214	2.21	ng	# 78

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

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