

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092215\
 Data File : BF081745.D
 Acq On : 22 Sep 2015 18:12
 Operator : UM/IZ
 Sample : G3753-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-2

Manual Integrations
 APPROVED

MMDadoda
 9/23/2015 4:33:44 PM

Quant Time: Sep 23 13:55:25 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 22 17:42:58 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.14	152	36078	20.00	ng	0.00
21) Naphthalene-d8	11.03	136	138035	20.00	ng	-0.01
38) Acenaphthene-d10	15.16	164	59368	20.00	ng	-0.01
63) Phenanthrene-d10	18.67	188	99177	20.00	ng	-0.01
75) Chrysene-d12	23.33	240	71240	20.00	ng	0.00
86) Perylene-d12	24.77	264	72934	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	268149	115.32	ng	0.02
7) Phenol-d6	7.58	99	371680	120.75	ng	0.00
23) Nitrobenzene-d5	9.44	82	247930	86.66	ng	-0.01
41) 2,4,6-Tribromophenol	17.07	330	61188	115.75	ng	-0.01
44) 2-Fluorobiphenyl	13.68	172	429860	92.79	ng	-0.01
78) Terphenyl-d14	22.05	244	258994	84.63	ng	0.00
Target Compounds						
49) Dimethylphthalate	14.73	163	93113	20.19	ng	Qvalue # 97
70) Phenanthrene	18.72	178	37618	6.82	ng	96
74) Fluoranthene	21.36	202	89039	14.92	ng	87
77) Pyrene	21.71	202	83127	15.75	ng	99
80) Benzo(a)anthracene	23.32	228	50414m	11.11	ng	
82) Chrysene	23.36	228	40201	9.89	ng	95
83) Bis(2-ethylhexyl)phthalate	23.44	149	7181	2.37	ng	# 95
85) Indeno(1,2,3-cd)pyrene	25.81	276	23835m	7.99	ng	
87) Benzo(b)fluoranthene	24.44	252	61571m	11.82	ng	
88) Benzo(k)fluoranthene	24.45	252	13966m	3.23	ng	
89) Benzo(a)pyrene	24.72	252	41006	9.29	ng	# 87
91) Benzo(a,h,i)perylene	26.11	276	24248	7.45	ng	# 72

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

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