

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF090815\
 Data File : BF081533.D
 Acq On : 8 Sep 2015 20:52
 Operator : UM/IZ
 Sample : G3555-10
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-10

Manual Integrations
 APPROVED

MMDadoda
 9/9/2015 2:58:52 PM

Quant Time: Sep 09 01:42:18 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 08 23:43:58 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.34	152	53596	20.00	ng	0.00
21) Naphthalene-d8	7.63	136	208254	20.00	ng	-0.01
38) Acenaphthene-d10	9.37	164	96911	20.00	ng	-0.01
63) Phenanthrene-d10	10.83	188	176034	20.00	ng	-0.01
75) Chrysene-d12	13.44	240	116766	20.00	ng	-0.01
86) Perylene-d12	14.75	264	109981	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.89	112	434897	125.90	ng	0.01
7) Phenol-d6	6.02	99	643975	140.84	ng	0.00
23) Nitrobenzene-d5	6.92	82	369429	85.59	ng	0.00
41) 2,4,6-Tribromophenol	10.16	330	110536	128.10	ng	0.00
44) 2-Fluorobiphenyl	8.72	172	568435	75.17	ng	0.00
78) Terphenyl-d14	12.42	244	459091	91.52	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.12	163	134411	17.85	ng	Qvalue # 97
70) Phenanthrene	10.86	178	32559	3.33	ng	97
74) Fluoranthene	12.03	202	73504	6.94	ng	87
77) Pyrene	12.25	202	59009	6.82	ng	96
80) Benzo(a)anthracene	13.44	228	31569m	4.25	ng	
82) Chrysene	13.47	228	24656m	3.70	ng	
85) Indeno(1,2,3-cd)pyrene	15.82	276	13199	2.70	ng	# 79
87) Benzo(b)fluoranthene	14.42	252	31127m	3.96	ng	
88) Benzo(k)fluoranthene	14.43	252	13745m	2.11	ng	
89) Benzo(a)pyrene	14.71	252	24152	3.63	ng	# 82
91) Benzo(g,h,i)perylene	16.14	276	11334	2.31	ng	# 71

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

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