

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091515\
 Data File : BF081636.D
 Acq On : 15 Sep 2015 19:12
 Operator : UM/IZ
 Sample : G3619-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 H1-H2-H3-H4-H5-COMP

Manual Integrations
 APPROVED

MMDadoda
 9/16/2015 1:24:03 PM

Quant Time: Sep 16 01:26:18 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 15 14:03:59 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.21	152	47760	20.00	ng	-0.01
21) Naphthalene-d8	11.11	136	189899	20.00	ng	-0.01
38) Acenaphthene-d10	15.25	164	85299	20.00	ng	-0.01
63) Phenanthrene-d10	18.76	188	159163	20.00	ng	0.00
75) Chrysene-d12	23.38	240	108979	20.00	ng	0.00
86) Perylene-d12	24.81	264	84761	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	294938	95.81	ng	0.02
7) Phenol-d6	7.63	99	418388	102.68	ng	-0.01
23) Nitrobenzene-d5	9.51	82	276551	70.26	ng	-0.02
41) 2,4,6-Tribromophenol	17.16	330	74749	98.42	ng	-0.01
44) 2-Fluorobiphenyl	13.76	172	463939	69.70	ng	-0.01
78) Terphenyl-d14	22.10	244	341841	73.02	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	14.81	163	119586	18.04	ng	# 97
70) Phenanthrene	18.80	178	92426	10.45	ng	98
74) Fluoranthene	21.43	202	144122	15.05	ng	86
77) Pyrene	21.77	202	152900	18.94	ng	99
80) Benzo(a)anthracene	23.37	228	64529	9.30	ng	97
82) Chrysene	23.41	228	61400	9.87	ng	94
85) Indeno(1,2,3-cd)pyrene	25.86	276	22157	4.85	ng	# 79
87) Benzo(b)fluoranthene	24.48	252	70377m	11.63	ng	
88) Benzo(k)fluoranthene	24.49	252	12228m	2.44	ng	
89) Benzo(a)pyrene	24.77	252	48804	9.52	ng	# 83
91) Benzo(g,h,i)perylene	26.16	276	19949	5.28	ng	# 73

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

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