

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091515\
 Data File : BF081629.D
 Acq On : 15 Sep 2015 14:48
 Operator : UM/IZ
 Sample : G3597-08 2X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GP-8(0-5)

Manual Integrations
 APPROVED

MMDadoda
 9/16/2015 1:23:20 PM

Quant Time: Sep 15 15:53:09 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	50498	20.00	ng	-0.01
21) Naphthalene-d8	11.11	136	201907	20.00	ng	-0.01
38) Acenaphthene-d10	15.25	164	95501	20.00	ng	-0.01
63) Phenanthrene-d10	18.74	188	177850	20.00	ng	-0.01
75) Chrysene-d12	23.37	240	110217	20.00	ng	-0.01
86) Perylene-d12	24.81	264	96531	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	241928	74.33	ng	0.01
7) Phenol-d6	7.62	99	332905	77.27	ng	-0.02
23) Nitrobenzene-d5	9.50	82	221168	52.85	ng	-0.03
41) 2,4,6-Tribromophenol	17.14	330	59664	70.16	ng	-0.02
44) 2-Fluorobiphenyl	13.75	172	359206	48.20	ng	-0.02
78) Terphenyl-d14	22.10	244	260444	55.01	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	14.81	163	87242	11.76	ng	# 96
70) Phenanthrene	18.80	178	108057	10.93	ng	98
71) Anthracene	18.93	178	24201	2.38	ng	96
74) Fluoranthene	21.42	202	192480	17.98	ng	91
77) Pyrene	21.77	202	176874	21.66	ng	96
80) Benzo(a)anthracene	23.37	228	76315	10.87	ng	96
82) Chrysene	23.41	228	70631	11.23	ng	95
85) Indeno(1,2,3-cd)pyrene	25.87	276	43023	9.32	ng	# 88
87) Benzo(b)fluoranthene	24.48	252	80688m	11.71	ng	
88) Benzo(k)fluoranthene	24.48	252	35617m	6.23	ng	
89) Benzo(a)pyrene	24.77	252	63602	10.89	ng	# 83
90) Dibenzo(a,h)anthracene	25.87	278	9227	2.04	ng	# 80
91) Benzo(g,h,i)perylene	26.16	276	44224	10.27	ng	# 75

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

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