

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080115\
 Data File : BF080801.D
 Acq On : 2 Aug 2015 1:42
 Operator : TP/UM
 Sample : G3125-01
 Misc :
 ALS Vial : 87 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WEST-ABUT-2(0-2)

Manual Integrations
 APPROVED

MMDadoda
 8/3/2015 3:42:32 PM

Quant Time: Aug 03 07:35:06 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF073015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 31 01:45:22 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	189680	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	654178	20.00	ng	-0.01
38) Acenaphthene-d10	9.89	164	286850	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	422345	20.00	ng	0.00
75) Chrysene-d12	14.00	240	256325	20.00	ng	0.00
86) Perylene-d12	15.43	264	366949	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	952893	86.18	ng	0.01
7) Phenol-d6	6.49	99	1376800	91.46	ng	0.00
23) Nitrobenzene-d5	7.41	82	769939	56.99	ng	-0.01
41) 2,4,6-Tribromophenol	10.67	330	253614	85.21	ng	-0.01
44) 2-Fluorobiphenyl	9.22	172	1580396	82.20	ng	0.00
78) Terphenyl-d14	12.96	244	943289	69.60	ng	0.00
Target Compounds						Qvalue
32) Benzoic acid	7.87	122	103557	14.53	ng	97
49) Dimethylphthalate	9.61	163	557607	27.58	ng	99
70) Phenanthrene	11.39	178	42722	2.01	ng	96
74) Fluoranthene	12.58	202	87499	4.25	ng	96
77) Pyrene	12.81	202	75939	3.82	ng	99
80) Benzo(a)anthracene	13.99	228	36281	2.35	ng	96
82) Chrysene	14.02	228	31323m	2.09	ng	
85) Indeno(1,2,3-cd)pyrene	16.80	276	29072	4.65	ng	# 88
87) Benzo(b)fluoranthene	15.01	252	71326m	3.25	ng	
89) Benzo(a)pyrene	15.36	252	38671	2.15	ng	# 78

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

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