

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080315\
 Data File : BF080835.D
 Acq On : 4 Aug 2015 10:46
 Operator : TP
 Sample : G3145-03
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-SF-03-003-B

Manual Integrations
 APPROVED

MMDadoda
 8/5/2015 2:29:33 PM

Quant Time: Aug 04 12:23:23 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF080315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 04 01:58:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	166788	20.00	ng	-0.01
21) Naphthalene-d8	8.11	136	551115	20.00	ng	-0.01
38) Acenaphthene-d10	9.87	164	220240	20.00	ng	0.00
63) Phenanthrene-d10	11.34	188	348738	20.00	ng	0.00
75) Chrysene-d12	13.97	240	360946	20.00	ng	-0.01
86) Perylene-d12	15.39	264	331333	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	1387631	138.71	ng	0.01
7) Phenol-d6	6.46	99	1747630	129.87	ng	0.00
23) Nitrobenzene-d5	7.40	82	1287018	115.24	ng	0.00
41) 2,4,6-Tribromophenol	10.66	330	292984	138.82	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	1654909	113.58	ng	0.00
78) Terphenyl-d14	12.93	244	1094273	60.42	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.58	163	195133	13.26	ng	99
70) Phenanthrene	11.37	178	53029	2.92	ng	97
74) Fluoranthene	12.56	202	123225	7.09	ng	94
77) Pyrene	12.79	202	125975	4.71	ng	96
80) Benzo(a)anthracene	13.96	228	89232	4.33	ng	95
82) Chrysene	14.00	228	66213m	3.31	ng	
87) Benzo(b)fluoranthene	14.99	252	85168m	4.39	ng	
89) Benzo(a)pyrene	15.33	252	59300	3.59	ng	# 92
91) Benzo(g,h,i)perylene	17.17	276	33621	2.01	ng	# 91

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

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