

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042716\
 Data File : BF086443.D
 Acq On : 28 Apr 2016 00:26
 Operator : UM/SJ
 Sample : H2713-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EUT01-B6B9B11B15B10-C

Manual Integrations
 APPROVED

sohil
 4/28/2016 1:40:21 PM

Quant Time: Apr 28 02:23:12 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 27 14:20:52 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	130629	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	471595	20.00	ng	-0.01
38) Acenaphthene-d10	9.85	164	174340	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	278842	20.00	ng	0.00
75) Chrysene-d12	13.96	240	306291	20.00	ng	0.00
86) Perylene-d12	15.37	264	239360	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	806270	106.46	ng	0.01
7) Phenol-d6	6.44	99	955654	90.41	ng	0.00
23) Nitrobenzene-d5	7.37	82	586100	66.99	ng	-0.01
41) 2,4,6-Tribromophenol	10.64	330	164318	89.37	ng	0.00
44) 2-Fluorobiphenyl	9.17	172	957678	99.64	ng	0.00
78) Terphenyl-d14	12.91	244	682203	49.07	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.56	163	138363	10.54	ng	100
70) Phenanthrene	11.34	178	50145	3.43	ng	97
74) Fluoranthene	12.53	202	132056	8.94	ng	96
77) Pyrene	12.76	202	130142	5.90	ng	98
80) Benzo(a)anthracene	13.95	228	94089	5.76	ng	95
82) Chrysene	13.99	228	54438m	3.07	ng	
85) Indeno(1,2,3-cd)pyrene	16.73	276	31195	2.37	ng	# 88
87) Benzo(b)fluoranthene	14.97	252	85973m	6.11	ng	
88) Benzo(k)fluoranthene	14.98	252	45417m	3.09	ng	
89) Benzo(a)pyrene	15.31	252	56161	4.22	ng	# 87
91) Benzo(g,h,i)perylene	17.13	276	29869	2.74	ng	98

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

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