

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050416\
 Data File : BF086704.D
 Acq On : 5 May 2016 4:16
 Operator : UM/SJ
 Sample : H2829-11
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EUT01-TD-B4(5-12)-C

Manual Integrations
 APPROVED

sohil
 5/5/2016 3:19:48 PM

Quant Time: May 05 13:07:57 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 04 15:12:53 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	157109	40.00	ng	0.00
21) Naphthalene-d8	8.03	136	545721	40.00	ng	0.00
38) Acenaphthene-d10	9.78	164	226164	40.00	ng	0.00
63) Phenanthrene-d10	11.25	188	345680m	40.00	ng	0.00
75) Chrysene-d12	13.89	240	312643	40.00	ng	0.01
86) Perylene-d12	15.28	264	213642	40.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	1056725	111.35	ng	0.01
7) Phenol-d6	6.38	99	1257441	98.36	ng	0.01
23) Nitrobenzene-d5	7.31	82	848335	79.88	ng	0.00
41) 2,4,6-Tribromophenol	10.57	330	224796	101.61	ng	0.00
44) 2-Fluorobiphenyl	9.10	172	1129574	81.87	ng	0.00
78) Terphenyl-d14	12.84	244	829786	53.64	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.49	163	57161	3.39	ng	98
70) Phenanthrene	11.28	178	79513	4.28	ng	99
74) Fluoranthene	12.47	202	124007	6.66	ng	94
77) Pyrene	12.69	202	128404	5.15	ng	98
80) Benzo(a)anthracene	13.88	228	51300	2.90	ng	94
82) Chrysene	13.92	228	48571m	2.57	ng	
87) Benzo(b)fluoranthene	14.89	252	54352m	4.23	ng	
89) Benzo(a)pyrene	15.22	252	34270	2.90	ng	# 79
91) Benzo(g,h,i)perylene	17.00	276	25824	2.17	ng	96

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

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