

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042016\
 Data File : BF086324.D
 Acq On : 20 Apr 2016 22:37
 Operator : UM/SJ
 Sample : H2601-03
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RODNEY-AVE-PS(0-2)B

Manual Integrations
 APPROVED

sohil
 4/21/2016 7:11:00 PM

Quant Time: Apr 21 01:49:33 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 20 15:45:26 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	255813	20.00	ng	-0.03
21) Naphthalene-d8	8.14	136	1001191	20.00	ng	-0.03
38) Acenaphthene-d10	9.90	164	386399	20.00	ng	-0.03
63) Phenanthrene-d10	11.38	188	571844	20.00	ng	-0.03
75) Chrysene-d12	14.02	240	424714	20.00	ng	-0.03
86) Perylene-d12	15.45	264	363945	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	1466019	99.99	ng	-0.02
7) Phenol-d6	6.50	99	2116051	110.29	ng	-0.02
23) Nitrobenzene-d5	7.42	82	1246337	71.29	ng	-0.03
41) 2,4,6-Tribromophenol	10.69	330	331263	109.99	ng	-0.03
44) 2-Fluorobiphenyl	9.23	172	1920216	98.29	ng	-0.03
78) Terphenyl-d14	12.97	244	1062900	59.57	ng	-0.03
Target Compounds						Qvalue
49) Dimethylphthalate	9.62	163	297293	10.11	ng	99
70) Phenanthrene	11.40	178	233441	8.61	ng	99
71) Anthracene	11.46	178	86172	2.97	ng	99
74) Fluoranthene	12.59	202	514035	20.46	ng	100
77) Pyrene	12.82	202	447128	14.19	ng	99
80) Benzo(a)anthracene	14.01	228	256921	11.48	ng	98
82) Chrysene	14.04	228	301087m	12.45	ng	
83) Bis(2-ethylhexyl)phthalate	14.01	149	58522	3.00	ng	99
85) Indeno(1,2,3-cd)pyrene	16.84	276	139156	5.60	ng	98
87) Benzo(b)fluoranthene	15.04	252	394432m	18.74	ng	
88) Benzo(k)fluoranthene	15.06	252	81497m	4.60	ng	
89) Benzo(a)pyrene	15.39	252	217066	11.19	ng	# 94
91) Benzo(a,h,i)perylene	17.27	276	147241	9.07	ng	# 94

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

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