

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042916\
 Data File : BF086498.D
 Acq On : 29 Apr 2016 18:03
 Operator : UM/SJ
 Sample : H2784-03
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP3(3A-3B-3C-3D)

Manual Integrations
 APPROVED

sohil
 4/30/2016 11:22:47 AM

Quant Time: Apr 30 01:24:10 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	6.77	152	162190	20.00	ng	-0.03
21) Naphthalene-d8	8.06	136	683770	20.00	ng	-0.03
38) Acenaphthene-d10	9.81	164	284287	20.00	ng	-0.03
63) Phenanthrene-d10	11.30	188	458542	20.00	ng	-0.02
75) Chrysene-d12	13.93	240	318575	20.00	ng	-0.03
86) Perylene-d12	15.33	264	263600	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	1138140	121.04	ng	-0.01
7) Phenol-d6	6.42	99	1752409	133.53	ng	-0.02
23) Nitrobenzene-d5	7.34	82	1056289	83.27	ng	-0.03
41) 2,4,6-Tribromophenol	10.60	330	308901	103.04	ng	-0.03
44) 2-Fluorobiphenyl	9.14	172	1568757	100.23	ng	-0.03
78) Terphenyl-d14	12.89	244	1047423	72.43	ng	-0.02
Target Compounds						
49) Dimethylphthalate	9.54	163	142779	6.67	ng	99
70) Phenanthrene	11.32	178	117748	4.89	ng	98
74) Fluoranthene	12.50	202	213935	8.81	ng	98
77) Pyrene	12.73	202	211885	9.23	ng	99
80) Benzo(a)anthracene	13.92	228	109697	6.46	ng	95
82) Chrysene	13.95	228	92515m	5.01	ng	
85) Indeno(1,2,3-cd)pyrene	16.67	276	64143	4.69	ng	# 92
87) Benzo(b)fluoranthene	14.93	252	119769m	7.72	ng	
88) Benzo(k)fluoranthene	14.95	252	67167m	4.15	ng	
89) Benzo(a)pyrene	15.28	252	92056	6.29	ng	# 84
91) Benzo(a,h,i)perylene	17.07	276	73627	6.13	ng	98

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

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