

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
 Data File : BF085588.D
 Acq On : 19 Mar 2016 22:19
 Operator : UM/SJ
 Sample : H1803-01RE
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-01RE

Manual Integrations
 APPROVED

apatel
 3/21/2016 5:09:27 PM

Quant Time: Mar 21 03:59:41 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 18 23:31:39 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.01	152	347456	20.00	ng	0.15
21) Naphthalene-d8	0.00	136	0	0.00	ng	-8.14
38) Acenaphthene-d10	0.00	164	0	0.00	ng	-9.90
63) Phenanthrene-d10	11.37	188	219	20.00	ng	-0.01
75) Chrysene-d12	14.01	240	1488	20.00	ng	0.00
86) Perylene-d12	15.43	264	1851	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	820968	39.69	ng	0.00
7) Phenol-d6	6.48	99	1031872	38.89	ng	-0.01
23) Nitrobenzene-d5	7.42	82	666314	0.00	ng	-0.01
41) 2,4,6-Tribromophenol	10.68	330	182720	0.00	ng	-0.01
44) 2-Fluorobiphenyl	9.21	172	955793	0.00	ng	-0.01
78) Terphenyl-d14	12.96	244	748157	12099.05	ng	0.00
Target Compounds						Qvalue
70) Phenanthrene	11.40	178	131791	11413.06	ng	99
71) Anthracene	11.44	178	37529	3099.58	ng	97
72) Carbazole	11.60	167	16773	1605.76	ng	99
74) Fluoranthene	12.58	202	209456	17323.35	ng	90
77) Pyrene	12.81	202	183434	1716.89	ng	99
80) Benzo(a)anthracene	13.99	228	101660	1176.84	ng	96
82) Chrysene	14.03	228	76035m	914.92	ng	
85) Indeno(1,2,3-cd)pyrene	16.84	276	49526	713.17	ng	92
87) Benzo(b)fluoranthene	15.03	252	130286m	1078.73	ng	
88) Benzo(k)fluoranthene	15.05	252	36066m	362.45	ng	
89) Benzo(a)pyrene	15.37	252	90986	888.61	ng	96
90) Dibenzo(a,h)anthracene	16.84	278	12177	142.59	ng	# 66
91) Benzo(g,h,i)perylene	17.25	276	50267	608.11	ng	92

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

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