

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062416\
 Data File : BF088334.D
 Acq On : 24 Jun 2016 16:27
 Operator : UM/SJ
 Sample : H3683-09DL 10X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-06-COMPD

Manual Integrations
 APPROVED

sohil
 6/27/2016 8:11:45 PM

Quant Time: Jun 24 18:00:14 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 22 14:55:14 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.58	152	77218m	20.00	ng	-0.02
21) Naphthalene-d8	7.87	136	312359	20.00	ng	-0.02
38) Acenaphthene-d10	9.61	164	147492	20.00	ng	-0.03
63) Phenanthrene-d10	11.10	188	277294	20.00	ng	-0.02
75) Chrysene-d12	13.73	240	170353	20.00	ng	-0.02
86) Perylene-d12	15.09	264	168411	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.19	112	66888	14.81	ng	0.01
7) Phenol-d6	6.25	99	99337	18.15	ng	-0.01
23) Nitrobenzene-d5	7.16	82	50235	9.39	ng	-0.01
41) 2,4,6-Tribromophenol	10.41	330	19820	12.73	ng	-0.02
44) 2-Fluorobiphenyl	8.95	172	122960	8.09	ng	-0.02
78) Terphenyl-d14	12.67	244	102369	13.67	ng	-0.02
Target Compounds						Qvalue
48) Acenaphthylene	9.48	152	66216	4.36	ng	98
70) Phenanthrene	11.12	178	60420	4.15	ng	98
71) Anthracene	11.17	178	109091	7.43	ng	97
74) Fluoranthene	12.30	202	635898	41.41	ng	98
77) Pyrene	12.53	202	586228	46.37	ng	100
80) Benzo(a)anthracene	13.71	228	440791	45.41	ng	# 79
82) Chrysene	13.75	228	209448m	22.93	ng	
85) Indeno(1,2,3-cd)pyrene	16.33	276	181201	21.35	ng	# 100
87) Benzo(b)fluoranthene	14.72	252	403346m	34.36	ng	
88) Benzo(k)fluoranthene	14.73	252	116236m	13.13	ng	
89) Benzo(a)pyrene	15.03	252	303780m	31.99	ng	
90) Dibenzo(a,h)anthracene	16.33	278	44476m	5.04	ng	
91) Benzo(g,h,i)perylene	16.70	276	151516	16.70	ng	97

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

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