

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080115\
 Data File : BF080788.D
 Acq On : 1 Aug 2015 19:01
 Operator : TP/UM
 Sample : G3119-01
 Misc :
 ALS Vial : 74 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BB-026CS-1(0-16FTBGS)

Manual Integrations
 APPROVED

MMDadoda
 8/3/2015 3:41:23 PM

Quant Time: Aug 03 06:59:50 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF073015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 31 01:45:22 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	90909	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	310698	20.00	ng	-0.01
38) Acenaphthene-d10	9.88	164	129883	20.00	ng	-0.01
63) Phenanthrene-d10	11.37	188	222561	20.00	ng	0.00
75) Chrysene-d12	14.00	240	187069	20.00	ng	0.00
86) Perylene-d12	15.41	264	299729	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	783166	147.79	ng	0.01
7) Phenol-d6	6.48	99	878250	121.73	ng	-0.01
23) Nitrobenzene-d5	7.41	82	636778	99.25	ng	-0.01
41) 2,4,6-Tribromophenol	10.67	330	192212	142.62	ng	-0.01
44) 2-Fluorobiphenyl	9.21	172	807364	92.74	ng	-0.01
78) Terphenyl-d14	12.95	244	525973	53.18	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.61	163	176928	19.33	ng	Qvalue # 96
70) Phenanthrene	11.39	178	28987	2.59	ng	99
74) Fluoranthene	12.57	202	61622	5.68	ng	99
77) Pyrene	12.80	202	66194	4.56	ng	98
80) Benzo(a)anthracene	13.99	228	39647	3.52	ng	93
82) Chrysene	14.02	228	34928m	3.20	ng	
85) Indeno(1,2,3-cd)pyrene	16.79	276	33374m	5.81	ng	
87) Benzo(b)fluoranthene	15.00	252	62712m	3.50	ng	
88) Benzo(k)fluoranthene	15.01	252	31659m	2.15	ng	
89) Benzo(a)pyrene	15.35	252	56670	3.86	ng	# 90
91) Benzo(g,h,i)perylene	17.20	276	34742	2.68	ng	95

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

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