

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030215\
 Data File : BF077536.D
 Acq On : 2 Mar 2015 21:46
 Operator : TP/IZ
 Sample : G1399-12
 Misc : S
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB6

Manual Integrations
 APPROVED

apatel
 3/3/2015 6:26:15 PM

Quant Time: Mar 03 04:34:05 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 02 23:46:07 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.26	152	87375	20.00	ng	0.00
21) Naphthalene-d8	8.84	136	351643	20.00	ng	0.00
38) Acenaphthene-d10	11.01	164	179707	20.00	ng	0.00
63) Phenanthrene-d10	12.85	188	365697	20.00	ng	0.00
75) Chrysene-d12	16.14	240	407268	20.00	ng	0.00
86) Perylene-d12	17.86	264	393178	20.00	ng	0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	414461	79.19	ng	0.00
7) Phenol-d6	6.82	99	505922	75.18	ng	0.00
23) Nitrobenzene-d5	7.95	82	300482	53.14	ng	0.00
41) 2,4,6-Tribromophenol	11.99	330	139756	80.77	ng	0.00
44) 2-Fluorobiphenyl	10.18	172	614647	53.33	ng	0.00
78) Terphenyl-d14	14.85	244	751916	43.16	ng	0.00
Target Compounds						Qvalue
48) Acenaphthylene	10.84	152	38210	2.26	ng	99
70) Phenanthrene	12.88	178	85212	4.02	ng	99
71) Anthracene	12.95	178	46580	2.21	ng	99
74) Fluoranthene	14.36	202	335908	14.51	ng	99
77) Pyrene	14.64	202	326292	13.10	ng	# 95
80) Benzo(a)anthracene	16.13	228	261324m	10.95	ng	
82) Chrysene	16.17	228	247240m	11.03	ng	
85) Indeno(1,2,3-cd)pyrene	19.31	276	201718m	7.89	ng	
87) Benzo(b)fluoranthene	17.43	252	397820m	17.40	ng	
88) Benzo(k)fluoranthene	17.45	252	142687m	6.37	ng	
89) Benzo(a)pyrene	17.80	252	301202	13.83	ng	# 95
90) Dibenzo(a,h)anthracene	19.33	278	49157	2.37	ng	# 89
91) Benzo(g,h,i)perylene	19.74	276	184581	8.81	ng	96

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

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