

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051515\
 Data File : BF079292.D
 Acq On : 16 May 2015 4:21
 Operator : TP/IZ
 Sample : G2215-09DL 4X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-38SDL

Manual Integrations
 APPROVED

apatel
 5/18/2015 3:41:31 PM

Quant Time: May 16 05:07:31 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat May 16 04:02:20 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.18	152	55954	20.00	ng	0.00
21) Naphthalene-d8	8.75	136	227030	20.00	ng	0.00
38) Acenaphthene-d10	10.92	164	104032	20.00	ng	0.00
63) Phenanthrene-d10	12.77	188	196825	20.00	ng	0.00
75) Chrysene-d12	16.05	240	199210	20.00	ng	-0.01
86) Perylene-d12	17.73	264	209292	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	77316m	23.79	ng	0.01
7) Phenol-d6	6.74	99	100829	23.47	ng	0.00
23) Nitrobenzene-d5	7.86	82	51401	13.01	ng	-0.01
41) 2,4,6-Tribromophenol	11.91	330	21260	18.89	ng	0.00
44) 2-Fluorobiphenyl	10.10	172	120220	16.10	ng	0.00
78) Terphenyl-d14	14.75	244	125202	15.05	ng	-0.01
Target Compounds						Qvalue
48) Acenaphthylene	10.75	152	22978	2.17	ng	98
70) Phenanthrene	12.79	178	220177	20.00	ng	99
71) Anthracene	12.86	178	54832	5.08	ng	99
74) Fluoranthene	14.27	202	451850	39.38	ng	92
77) Pyrene	14.55	202	407379	35.49	ng	99
80) Benzo(a)anthracene	16.03	228	233465	21.40	ng	97
82) Chrysene	16.08	228	176141	16.79	ng	98
85) Indeno(1,2,3-cd)pyrene	19.12	276	154236	11.54	ng	# 94
87) Benzo(b)fluoranthene	17.30	252	264959m	23.13	ng	
88) Benzo(k)fluoranthene	17.31	252	113513m	10.24	ng	
89) Benzo(a)pyrene	17.67	252	209317	19.84	ng	98
90) Dibenzo(a,h)anthracene	19.13	278	31827	2.84	ng	# 89
91) Benzo(a,h,i)perylene	19.52	276	152833	13.60	ng	# 95

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

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