

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061615\
 Data File : BF080014.D
 Acq On : 17 Jun 2015 5:29
 Operator : TP/IZ
 Sample : G2651-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NC-TS-001

Manual Integrations
 APPROVED

apatel
 6/17/2015 5:13:03 PM

Quant Time: Jun 17 06:59:24 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 15 15:44:31 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.35	152	51810	20.00	ng	0.00
21) Naphthalene-d8	8.64	136	207937	20.00	ng	0.00
38) Acenaphthene-d10	10.41	164	106800	20.00	ng	0.00
63) Phenanthrene-d10	11.92	188	216526	20.00	ng	0.00
75) Chrysene-d12	14.93	240	247218	20.00	ng	0.02
86) Perylene-d12	16.85	264	246005	20.00	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.97	112	283079	100.19	ng	0.01
7) Phenol-d6	6.95	99	399064	103.67	ng	0.00
23) Nitrobenzene-d5	7.91	82	216683	68.75	ng	0.00
41) 2,4,6-Tribromophenol	11.21	330	98799	92.72	ng	0.00
44) 2-Fluorobiphenyl	9.72	172	462261	60.61	ng	0.00
78) Terphenyl-d14	13.64	244	583037	54.37	ng	0.01
Target Compounds						Qvalue
49) Dimethylphthalate	10.10	163	45013	5.26	ng	98
74) Fluoranthene	13.23	202	66427	4.66	ng	96
77) Pyrene	13.48	202	61445	3.96	ng	98
80) Benzo(a)anthracene	14.92	228	38945	2.48	ng	93
82) Chrysene	14.96	228	49235	2.39	ng	97
85) Indeno(1,2,3-cd)pyrene	18.72	276	39302	2.39	ng	# 94
87) Benzo(b)fluoranthene	16.29	252	68834m	4.57	ng	
89) Benzo(a)pyrene	16.76	252	44605	3.11	ng	# 94
91) Benzo(g,h,i)perylene	19.28	276	45167	3.10	ng	96

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

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