

Data Path : Z:\HPCHEM1\BNA F\DATA\BF061115\
 Data File : BF079915.D
 Acq On : 12 Jun 2015 2:25
 Operator : TP/IZ
 Sample : G2449-09
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FM6

Manual Integrations
 APPROVED

apatel
 6/13/2015 8:44:41 AM

Quant Time: Jun 12 05:42:29 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF061115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 12 00:59:58 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	66992	20.00	ng	0.00
21) Naphthalene-d8	8.68	136	272429	20.00	ng	0.00
38) Acenaphthene-d10	10.46	164	149586	20.00	ng	0.00
63) Phenanthrene-d10	11.97	188	330857	20.00	ng	-0.01
75) Chrysene-d12	14.83	240	362165	20.00	ng	-0.16
86) Perylene-d12	16.69	264	356821	20.00	ng	-0.24
System Monitoring Compounds						
5) 2-Fluorophenol	6.01	112	413450	103.41	ng	0.01
7) Phenol-d6	6.99	99	592731	113.32	ng	0.00
23) Nitrobenzene-d5	7.95	82	296250	64.68	ng	0.00
41) 2,4,6-Tribromophenol	11.26	330	151597	113.55	ng	0.00
44) 2-Fluorobiphenyl	9.76	172	675340	63.63	ng	0.00
78) Terphenyl-d14	13.61	244	866567	54.37	ng	-0.08
Target Compounds						Qvalue
49) Dimethylphthalate	10.15	163	51064	4.33	ng	98
70) Phenanthrene	11.99	178	418676	20.96	ng	99
71) Anthracene	12.05	178	104065	5.39	ng	98
74) Fluoranthene	13.22	202	564513	26.89	ng	99
77) Pyrene	13.47	202	498656	22.24	ng	100
80) Benzo(a)anthracene	14.82	228	289795	13.47	ng	98
82) Chrysene	14.87	228	213186	10.57	ng	97
85) Indeno(1,2,3-cd)pyrene	18.57	276	137851	6.40	ng	96
87) Benzo(b)fluoranthene	16.14	252	285434m	13.45	ng	
88) Benzo(k)fluoranthene	16.17	252	97491m	4.42	ng	
89) Benzo(a)pyrene	16.61	252	207710	10.36	ng	96
90) Dibenzo(a,h)anthracene	18.59	278	45806	2.47	ng	97
91) Benzo(a,h,i)perylene	19.15	276	129960	6.63	ng	96

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

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