

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030215\
 Data File : BF077535.D
 Acq On : 2 Mar 2015 21:18
 Operator : TP/IZ
 Sample : G1399-09
 Misc : S
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB3

Manual Integrations
 APPROVED

apatel
 3/3/2015 6:26:18 PM

Quant Time: Mar 03 04:28:09 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	89111	20.00	ng	0.00
21) Naphthalene-d8	8.84	136	361131	20.00	ng	0.00
38) Acenaphthene-d10	11.01	164	184770	20.00	ng	0.00
63) Phenanthrene-d10	12.85	188	360556	20.00	ng	0.00
75) Chrysene-d12	16.15	240	398267	20.00	ng	0.01
86) Perylene-d12	17.89	264	373945	20.00	ng	0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	447117	83.76	ng	0.00
7) Phenol-d6	6.82	99	522074	76.07	ng	0.00
23) Nitrobenzene-d5	7.95	82	310663	53.49	ng	0.00
41) 2,4,6-Tribromophenol	11.99	330	143714	80.78	ng	0.00
44) 2-Fluorobiphenyl	10.18	172	648984	54.77	ng	0.00
78) Terphenyl-d14	14.85	244	759958	44.61	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	10.69	163	30662	2.36	ng	# 95
70) Phenanthrene	12.88	178	104714	5.01	ng	98
74) Fluoranthene	14.36	202	231446	10.14	ng	100
77) Pyrene	14.64	202	244407	10.03	ng	96
80) Benzo(a)anthracene	16.14	228	148550	6.36	ng	95
82) Chrysene	16.17	228	112517	5.13	ng	97
85) Indeno(1,2,3-cd)pyrene	19.35	276	76638m	3.07	ng	
87) Benzo(b)fluoranthene	17.45	252	175366m	8.06	ng	
88) Benzo(k)fluoranthene	17.47	252	49373m	2.32	ng	
89) Benzo(a)pyrene	17.82	252	121957	5.89	ng	98
91) Benzo(g,h,i)perylene	19.77	276	84562	4.24	ng	94

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

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