

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
 Data File : BF099202.D
 Acq On : 7 Oct 2017 18:56
 Operator : SJ/JU
 Sample : I5586-07 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 G47A-2.5-5

Manual Integrations
 APPROVED

Sohil
 10/9/2017 6:34:21 PM

Quant Time: Oct 08 04:51:11 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Oct 08 03:54:40 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	120159	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	446390	20.00	ng	0.00
38) Acenaphthene-d10	9.90	164	172649	20.00	ng	0.00
63) Phenanthrene-d10	11.39	188	267535	20.00	ng	0.00
75) Chrysene-d12	14.03	240	218787	20.00	ng	0.00
86) Perylene-d12	15.51	264	171869	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	147774	19.58	ng	0.00
7) Phenol-d6	6.49	99	174044	18.69	ng	0.00
23) Nitrobenzene-d5	7.42	82	91184	13.10	ng	-0.01
41) 2,4,6-Tribromophenol	10.69	330	21795	15.43	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	172131	16.64	ng	0.00
78) Terphenyl-d14	12.97	244	119095	12.38	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.61	163	28712	2.203	ng	98
74) Fluoranthene	12.60	202	162196	11.185	ng	98
77) Pyrene	12.83	202	145564	8.725	ng	98
80) Benzo(a)anthracene	14.02	228	51188	3.864	ng	95
82) Chrysene	14.06	228	43796	3.472	ng	96
87) Benzo(b)fluoranthene	15.07	252	49084m	4.645	ng	
89) Benzo(a)pyrene	15.44	252	22268	2.296	ng	# 79

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

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