

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080917\
 Data File : BF097541.D
 Acq On : 10 Aug 2017 3:07
 Operator : SJ/JU
 Sample : I4677-09
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-F10-BASE

Manual Integrations
 APPROVED

Sohil
 8/10/2017 6:08:00 PM

Quant Time: Aug 10 07:35:09 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 07 16:02:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.40	152	104647	20.00	ng	-0.01
21) Naphthalene-d8	7.68	136	440105	20.00	ng	-0.02
38) Acenaphthene-d10	9.42	164	179954	20.00	ng	-0.02
63) Phenanthrene-d10	10.89	188	257464	20.00	ng	-0.02
75) Chrysene-d12	13.52	240	203103	20.00	ng	-0.02
86) Perylene-d12	14.86	264	188038	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	4.99	112	832186	119.88	ng	-0.01
7) Phenol-d6	6.07	99	936853	118.22	ng	-0.01
23) Nitrobenzene-d5	6.97	82	637675	85.00	ng	-0.02
41) 2,4,6-Tribromophenol	10.21	330	147479	103.73	ng	-0.02
44) 2-Fluorobiphenyl	8.76	172	930114	87.72	ng	-0.02
78) Terphenyl-d14	12.48	244	517106	51.91	ng	-0.02
Target Compounds						
49) Dimethylphthalate	9.16	163	162112	11.863	ng	98
70) Phenanthrene	10.92	178	123197	8.931	ng	99
74) Fluoranthene	12.10	202	126082	9.329	ng	97
77) Pyrene	12.32	202	116117	6.983	ng	98
80) Benzo(a)anthracene	13.51	228	43704	3.326	ng	97
82) Chrysene	13.54	228	38628	3.010	ng	97
87) Benzo(b)fluoranthene	14.52	252	44061m	3.898	ng	
89) Benzo(a)pyrene	14.81	252	35723	3.453	ng	98

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

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