

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090117\
 Data File : BF098223.D
 Acq On : 1 Sep 2017 13:40
 Operator : SJ/JU
 Sample : I4987-04MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-06-COMPMSD

Manual Integrations
 APPROVED

mohammad
 9/5/2017 5:22:08 PM

Quant Time: Sep 01 16:38:19 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 01 16:25:25 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	233430	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	894782	20.00	ng	0.00
38) Acenaphthene-d10	9.76	164	376451	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	594592	20.00	ng	0.00
75) Chrysene-d12	13.87	240	453242	20.00	ng	0.00
86) Perylene-d12	15.29	264	428119	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	933939	60.86	ng	0.01
7) Phenol-d6	6.37	99	1288415	61.79	ng	0.00
23) Nitrobenzene-d5	7.29	82	809535	49.15	ng	0.00
41) 2,4,6-Tribromophenol	10.55	330	194086	59.42	ng	0.00
44) 2-Fluorobiphenyl	9.08	172	1037476	40.49	ng	0.00
78) Terphenyl-d14	12.82	244	641507	29.52	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.47	163	195826	7.116	ng	98
70) Phenanthrene	11.26	178	112434	3.549	ng	99
74) Fluoranthene	12.44	202	139546	4.220	ng	99
77) Pyrene	12.67	202	135956	3.668	ng	98
80) Benzo(a)anthracene	13.86	228	65643	2.416	ng	98
82) Chrysene	13.89	228	58708	2.173	ng	98
87) Benzo(b)fluoranthene	14.88	252	67680m	2.508	ng	
89) Benzo(a)pyrene	15.22	252	49109	2.056	ng	98

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

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