

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030917\
 Data File : BM009159.D
 Acq On : 09 Mar 2017 23:21
 Operator : SJ/MA
 Sample : I2062-17DL 5X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AK3DL

Manual Integrations
 APPROVED

mohammad
 3/10/2017 6:01:41 PM

Quant Time: Mar 10 03:54:56 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM030317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 10 03:16:46 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	302	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	905	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	504	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	1139	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	1275	0.40	ng/ul	0.00
20) Perylene-d12	23.78	264	1312	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	74	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	195	0.05	ng/ul	0.00
Target Compounds						
12) Phenanthrene	17.30	178	491	0.14	ng/ul#	71
15) Fluoranthene	19.32	202	1654	0.35	ng/ul	98
17) Pyrene	19.68	202	1585	0.36	ng/ul	93
18) Benzo(a)anthracene	21.42	228	1331	0.31	ng/ul	93
19) Chrysene	21.47	228	1556m	0.39	ng/ul	
21) Benzo(b)fluoranthene	23.07	252	2695m	0.52	ng/ul	
22) Benzo(k)fluoranthene	23.11	252	667m	0.15	ng/ul	
23) Benzo(a)pyrene	23.67	252	1778m	0.38	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.15	276	1203	0.20	ng/ul#	88
25) Dibenzo(a,h)anthracene	26.15	278	291	0.06	ng/ul#	16
26) Benzo(g,h,i)perylene	26.89	276	1109	0.22	ng/ul#	91

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

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