

Data Path : Z:\HPCHEM1\BNA M\DATA\BM070716\
 Data File : BM006302.D
 Acq On : 07 Jul 2016 03:29
 Operator : UM/SJ
 Sample : H3811-18
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4TW6

Manual Integrations

umangi
 7/7/2016 6:06:39 PM

Quant Time: Jul 07 12:37:32 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM070716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 07 05:51:09 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	5001	0.40	ng/ul	0.00
4) Naphthalene-d8	10.43	136	19625	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.30	164	11189	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.04	188	22126m	0.40	ng/ul	0.00
18) Chrysene-d12	21.24	240	14418	0.40	ng/ul	0.00
22) Perylene-d12	23.45	264	12376m	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.17	96	11789	5.58	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.02	152	9582	0.42	ng/ul	0.00
16) Fluoranthene-d10	19.08	212	17467	0.37	ng/ul	0.00
Target Compounds				Ovalue		
9) Acenaphthylene	14.01	152	3148	0.08	ng/ul#	95
14) Phenanthrene	17.07	178	21340m	0.35	ng/ul	
15) Anthracene	17.18	178	2808	0.05	ng/ul#	79
17) Fluoranthene	19.10	202	39858m	0.59	ng/ul	
19) Pyrene	19.48	202	36444	0.60	ng/ul#	91
20) Benzo(a)anthracene	21.22	228	11182	0.24	ng/ul#	84
21) Chrysene	21.28	228	14643	0.32	ng/ul	95
23) Benzo(b)fluoranthene	22.79	252	17910	0.40	ng/ul#	78
24) Benzo(k)fluoranthene	22.81	252	7174m	0.17	ng/ul	
25) Benzo(a)pyrene	23.35	252	15923m	0.39	ng/ul	
26) Indeno(1,2,3-cd)pyrene	25.66	276	9441m	0.21	ng/ul	
27) Dibenzo(a,h)anthracene	25.67	278	2375m	0.07	ng/ul	
28) Benzo(g,h,i)perylene	26.34	276	9665m	0.25	ng/ul	

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

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