

Data Path : Z:\HPCHEM1\BNA M\DATA\BM070716\
 Data File : BM006304.D
 Acq On : 07 Jul 2016 04:41
 Operator : UM/SJ
 Sample : H3811-20
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4TW8

Manual Integrations

umangi
 7/7/2016 6:06:14 PM

Quant Time: Jul 07 12:47:36 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	5131	0.40	ng/ul	0.00
4) Naphthalene-d8	10.43	136	20372	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.30	164	11821	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.04	188	24027m	0.40	ng/ul	0.00
18) Chrysene-d12	21.24	240	15442	0.40	ng/ul	0.00
22) Perylene-d12	23.45	264	11981m	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.17	96	11811	5.45	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.02	152	9944	0.42	ng/ul	0.00
16) Fluoranthene-d10	19.07	212	19700	0.39	ng/ul	-0.01
Target Compounds				Ovalue		
9) Acenaphthylene	14.01	152	3884	0.10	ng/ul#	94
14) Phenanthrene	17.07	178	26491m	0.40	ng/ul	
15) Anthracene	17.18	178	3308	0.05	ng/ul#	84
17) Fluoranthene	19.11	202	55569m	0.76	ng/ul	
19) Pyrene	19.47	202	48853	0.76	ng/ul	96
20) Benzo(a)anthracene	21.22	228	12931	0.26	ng/ul#	89
21) Chrysene	21.27	228	19899	0.41	ng/ul	95
23) Benzo(b)fluoranthene	22.78	252	25940m	0.60	ng/ul	
24) Benzo(k)fluoranthene	22.83	252	7180m	0.18	ng/ul	
25) Benzo(a)pyrene	23.35	252	13390m	0.34	ng/ul	
26) Indeno(1,2,3-cd)pyrene	25.66	276	11275m	0.25	ng/ul	
27) Dibenzo(a,h)anthracene	25.67	278	2678m	0.08	ng/ul	
28) Benzo(g,h,i)perylene	26.34	276	11331m	0.30	ng/ul	

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

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