

Data Path : Z:\HPCHEM1\BNA M\DATA\BM070716\
 Data File : BM006305.D
 Acq On : 07 Jul 2016 05:17
 Operator : UM/SJ
 Sample : H3811-21
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4TW9

Manual Integrations

umangi
 7/7/2016 6:06:12 PM

Quant Time: Jul 07 13:01:58 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	5140	0.40	ng/ul	0.00
4) Naphthalene-d8	10.43	136	20686	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.30	164	12040	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.04	188	26139m	0.40	ng/ul	0.00
18) Chrysene-d12	21.24	240	18509	0.40	ng/ul	0.00
22) Perylene-d12	23.45	264	12514m	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.17	96	10811	4.98	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.02	152	9204	0.38	ng/ul	0.00
16) Fluoranthene-d10	19.08	212	20201	0.37	ng/ul	0.00
Target Compounds				Ovalue		
9) Acenaphthylene	14.01	152	3909	0.10	ng/ul#	94
14) Phenanthrene	17.07	178	25683m	0.36	ng/ul	
15) Anthracene	17.18	178	3324	0.05	ng/ul#	85
17) Fluoranthene	19.11	202	51048m	0.64	ng/ul	
19) Pyrene	19.47	202	47751	0.62	ng/ul	97
20) Benzo(a)anthracene	21.22	228	14258	0.24	ng/ul#	90
21) Chrysene	21.27	228	19067	0.32	ng/ul	94
23) Benzo(b)fluoranthene	22.78	252	22124m	0.49	ng/ul	
24) Benzo(k)fluoranthene	22.83	252	6478m	0.16	ng/ul	
25) Benzo(a)pyrene	23.35	252	12930m	0.31	ng/ul	
26) Indeno(1,2,3-cd)pyrene	25.66	276	9488m	0.21	ng/ul	
27) Dibenzo(a,h)anthracene	25.67	278	2579m	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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