

Data Path : Z:\HPCHEM1\BNA M\DATA\BM070716\
 Data File : BM006289.D
 Acq On : 06 Jul 2016 19:02
 Operator : UM/SJ
 Sample : H3811-03
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4TV3

Manual Integrations

umangi
 7/7/2016 6:04:23 PM

Quant Time: Jul 07 10:56:10 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	4264	0.40	ng/ul	0.00
4) Naphthalene-d8	10.43	136	17021	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.30	164	9942	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.04	188	20883m	0.40	ng/ul	0.00
18) Chrysene-d12	21.24	240	14362	0.40	ng/ul	0.00
22) Perylene-d12	23.45	264	10954m	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.17	96	9823	5.45	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.02	152	7564	0.38	ng/ul	0.00
16) Fluoranthene-d10	19.07	212	15976	0.36	ng/ul	-0.01
Target Compounds				Ovalue		
5) Naphthalene	10.48	128	1285	0.03	ng/ul#	77
7) 2-Methylnaphthalene	12.09	142	779	0.03	ng/ul	99
9) Acenaphthylene	14.01	152	6477	0.19	ng/ul#	94
14) Phenanthrene	17.07	178	16526m	0.29	ng/ul	
15) Anthracene	17.16	178	6447	0.12	ng/ul#	86
17) Fluoranthene	19.11	202	95179m	1.49	ng/ul	
19) Pyrene	19.47	202	100530	1.67	ng/ul	96
20) Benzo(a)anthracene	21.22	228	38475	0.83	ng/ul#	92
21) Chrysene	21.27	228	44109	0.97	ng/ul	94
23) Benzo(b)fluoranthene	22.78	252	69102m	1.76	ng/ul	
24) Benzo(k)fluoranthene	22.83	252	17473m	0.48	ng/ul	
25) Benzo(a)pyrene	23.35	252	33625m	0.92	ng/ul	
26) Indeno(1,2,3-cd)pyrene	25.66	276	24902m	0.61	ng/ul	
27) Dibenzo(a,h)anthracene	25.67	278	6792m	0.21	ng/ul	
28) Benzo(a,h,i)perylene	26.34	276	21668m	0.63	ng/ul	

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

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