

Data Path : Z:\HPCHEM1\BNA M\DATA\BM070716\
 Data File : BM006287.D
 Acq On : 06 Jul 2016 17:49
 Operator : UM/SJ
 Sample : H3811-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4TV1

Manual Integrations

umangi
 7/7/2016 6:04:30 PM

Quant Time: Jul 07 10:44:44 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	3828	0.40	ng/ul	0.00
4) Naphthalene-d8	10.43	136	15280	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.30	164	9051	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.04	188	19146m	0.40	ng/ul	0.00
18) Chrysene-d12	21.24	240	12795	0.40	ng/ul	0.00
22) Perylene-d12	23.45	264	9580m	0.40	ng/ul	0.00

System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.17	96	8749	5.41	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.02	152	6675	0.37	ng/ul	0.00
16) Fluoranthene-d10	19.07	212	13542	0.34	ng/ul	-0.01

Target Compounds				Ovalue		
5) Naphthalene	10.48	128	1209	0.03	ng/ul#	80
7) 2-Methylnaphthalene	12.09	142	679	0.03	ng/ul	98
9) Acenaphthylene	14.01	152	1119	0.04	ng/ul#	88
17) Fluoranthene	19.11	202	21466	0.37	ng/ul	95
19) Pyrene	19.47	202	20858	0.39	ng/ul	96
20) Benzo(a)anthracene	21.22	228	7617	0.18	ng/ul#	92
21) Chrysene	21.27	228	9786	0.24	ng/ul	95
23) Benzo(b)fluoranthene	22.78	252	13087m	0.38	ng/ul	
24) Benzo(k)fluoranthene	22.83	252	3342m	0.10	ng/ul	
25) Benzo(a)pyrene	23.35	252	7006m	0.22	ng/ul	
26) Indeno(1,2,3-cd)pyrene	25.66	276	5896m	0.17	ng/ul	
28) Benzo(g,h,i)perylene	26.34	276	5737m	0.19	ng/ul	

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

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