

Data Path : Z:\HPCHEM1\BNA M\DATA\BM070716\
 Data File : BM006306.D
 Acq On : 07 Jul 2016 05:53
 Operator : UM/SJ
 Sample : H3811-22
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4TX0

Manual Integrations

umangi
 7/7/2016 6:06:09 PM

Quant Time: Jul 07 13:08:20 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	4912	0.40	ng/ul	0.00
4) Naphthalene-d8	10.43	136	19875	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.30	164	11154	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.04	188	23166m	0.40	ng/ul	0.00
18) Chrysene-d12	21.25	240	14371	0.40	ng/ul	0.00
22) Perylene-d12	23.45	264	11624m	0.40	ng/ul	0.00

System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.17	96	8897	4.29	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.02	152	7252	0.31	ng/ul	0.00
16) Fluoranthene-d10	19.08	212	14383	0.29	ng/ul	0.00

Target Compounds				Ovalue		
9) Acenaphthylene	14.01	152	1643	0.04	ng/ul#	36
14) Phenanthrene	17.07	178	9526m	0.15	ng/ul	
15) Anthracene	17.16	178	1473	0.02	ng/ul#	31
17) Fluoranthene	19.10	202	20601	0.29	ng/ul	94
19) Pyrene	19.48	202	19386	0.32	ng/ul#	92
20) Benzo(a)anthracene	21.22	228	5455	0.12	ng/ul#	81
21) Chrysene	21.28	228	7496	0.16	ng/ul	92
23) Benzo(b)fluoranthene	22.79	252	10030m	0.24	ng/ul	
24) Benzo(k)fluoranthene	22.83	252	2561m	0.07	ng/ul	
25) Benzo(a)pyrene	23.35	252	5434m	0.14	ng/ul	
26) Indeno(1,2,3-cd)pyrene	25.66	276	4656m	0.11	ng/ul	
28) Benzo(g,h,i)perylene	26.34	276	4522m	0.12	ng/ul	

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

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