

Data Path : Z:\HPCHEM1\BNA M\DATA\BM070716\
 Data File : BM006295.D
 Acq On : 06 Jul 2016 22:40
 Operator : UM/SJ
 Sample : H3811-09
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4TV7

Manual Integrations

umangi
 7/7/2016 6:04:59 PM

Quant Time: Jul 07 11:24:41 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	4439	0.40	ng/ul	0.00
4) Naphthalene-d8	10.43	136	17863	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.30	164	10317	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.04	188	23102m	0.40	ng/ul	0.00
18) Chrysene-d12	21.24	240	16478	0.40	ng/ul	0.00
22) Perylene-d12	23.45	264	11677m	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.17	96	6894	3.67	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.02	152	5805	0.28	ng/ul	0.00
16) Fluoranthene-d10	19.07	212	13231	0.27	ng/ul	-0.01
Target Compounds				Ovalue		
14) Phenanthrene	17.07	178	6491m	0.10	ng/ul	
15) Anthracene	17.18	178	1495	0.03	ng/ul#	68
17) Fluoranthene	19.11	202	15420	0.22	ng/ul	93
19) Pyrene	19.47	202	14204	0.21	ng/ul#	93
20) Benzo(a)anthracene	21.22	228	5547	0.10	ng/ul#	91
21) Chrysene	21.27	228	6541	0.13	ng/ul	93
23) Benzo(b)fluoranthene	22.78	252	8151m	0.19	ng/ul	
24) Benzo(k)fluoranthene	22.81	252	3222m	0.08	ng/ul	
25) Benzo(a)pyrene	23.35	252	4376m	0.11	ng/ul	
26) Indeno(1,2,3-cd)pyrene	25.66	276	3760m	0.09	ng/ul	
28) Benzo(g,h,i)perylene	26.34	276	3477m	0.10	ng/ul	

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

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