

Data Path : Z:\HPCHEM1\BNA M\DATA\BM070716\
 Data File : BM006301.D
 Acq On : 07 Jul 2016 02:53
 Operator : UM/SJ
 Sample : H3811-17
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4TW5

Manual Integrations

umangi
 7/7/2016 6:06:20 PM

Quant Time: Jul 07 11:58:45 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	4835	0.40	ng/ul	0.00
4) Naphthalene-d8	10.43	136	19248	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.30	164	10800	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.04	188	22149m	0.40	ng/ul	0.00
18) Chrysene-d12	21.24	240	14457	0.40	ng/ul	0.00
22) Perylene-d12	23.45	264	12097m	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.17	96	6748	3.30	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.02	152	6177	0.27	ng/ul	0.00
16) Fluoranthene-d10	19.07	212	12610	0.27	ng/ul	-0.01
Target Compounds				Ovalue		
5) Naphthalene	10.48	128	3414	0.08	ng/ul#	92
7) 2-Methylnaphthalene	12.09	142	1734	0.06	ng/ul	100
9) Acenaphthylene	14.01	152	12264	0.34	ng/ul#	93
10) Acenaphthene	14.35	153	1144	0.03	ng/ul#	87
11) Fluorene	15.34	166	3331	0.08	ng/ul#	79
14) Phenanthrene	17.07	178	59635m	0.97	ng/ul	
15) Anthracene	17.16	178	13705	0.24	ng/ul#	92
17) Fluoranthene	19.11	202	124829m	1.85	ng/ul	
19) Pyrene	19.47	202	141022	2.33	ng/ul	97
20) Benzo(a)anthracene	21.22	228	52292	1.12	ng/ul#	91
21) Chrysene	21.27	228	67313	1.47	ng/ul	95
23) Benzo(b)fluoranthene	22.78	252	70217m	1.62	ng/ul	
24) Benzo(k)fluoranthene	22.83	252	17766m	0.44	ng/ul	
25) Benzo(a)pyrene	23.35	252	46513m	1.15	ng/ul	
26) Indeno(1,2,3-cd)pyrene	25.66	276	30710m	0.69	ng/ul	
27) Dibenzo(a,h)anthracene	25.67	278	8898m	0.25	ng/ul	
28) Benzo(g,h,i)perylene	26.34	276	25508m	0.67	ng/ul	

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

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