

Data Path : Z:\HPCHEM1\BNA M\DATA\BM070716\
 Data File : BM006296.D
 Acq On : 06 Jul 2016 23:16
 Operator : UM/SJ
 Sample : H3811-10
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4TV8

Manual Integrations

umangi
 7/7/2016 6:05:37 PM

Quant Time: Jul 07 11:29:24 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	4458	0.40	ng/ul	0.00
4) Naphthalene-d8	10.43	136	17803	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.30	164	10183	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.04	188	22610m	0.40	ng/ul	0.00
18) Chrysene-d12	21.24	240	15479	0.40	ng/ul	0.00
22) Perylene-d12	23.45	264	12090m	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.17	96	6564	3.48	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.02	152	5322	0.26	ng/ul	0.00
16) Fluoranthene-d10	19.07	212	11866	0.25	ng/ul	-0.01
Target Compounds				Ovalue		
5) Naphthalene	10.48	128	3384	0.08	ng/ul#	91
7) 2-Methylnaphthalene	12.09	142	1645	0.06	ng/ul	99
9) Acenaphthylene	14.01	152	6639	0.19	ng/ul#	94
10) Acenaphthene	14.35	153	2964	0.09	ng/ul	89
11) Fluorene	15.34	166	4024	0.11	ng/ul#	83
14) Phenanthrene	17.07	178	75311m	1.20	ng/ul	
15) Anthracene	17.16	178	16541	0.29	ng/ul#	93
17) Fluoranthene	19.11	202	140736m	2.04	ng/ul	
19) Pyrene	19.47	202	137192	2.12	ng/ul	96
20) Benzo(a)anthracene	21.22	228	65188	1.30	ng/ul#	93
21) Chrysene	21.27	228	66957	1.36	ng/ul	96
23) Benzo(b)fluoranthene	22.78	252	77964m	1.80	ng/ul	
24) Benzo(k)fluoranthene	22.83	252	19250m	0.48	ng/ul	
25) Benzo(a)pyrene	23.35	252	52780m	1.31	ng/ul	
26) Indeno(1,2,3-cd)pyrene	25.66	276	35395m	0.79	ng/ul	
27) Dibenzo(a,h)anthracene	25.66	278	9218m	0.26	ng/ul	
28) Benzo(g,h,i)perylene	26.34	276	34691m	0.92	ng/ul	

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

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