

Data Path : Z:\HPCHEM1\BNA M\DATA\BM070716\
 Data File : BM006298.D
 Acq On : 07 Jul 2016 00:28
 Operator : UM/SJ
 Sample : H3811-12
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4TW0

Manual Integrations

umangi
 7/7/2016 6:05:31 PM

Quant Time: Jul 07 11:38:48 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	4690	0.40	ng/ul	0.00
4) Naphthalene-d8	10.43	136	18532	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.30	164	10453	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.04	188	21200m	0.40	ng/ul	0.00
18) Chrysene-d12	21.24	240	13859m	0.40	ng/ul	0.00
22) Perylene-d12	23.44	264	13636	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.17	96	7071	3.57	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.02	152	5658	0.26	ng/ul	0.00
16) Fluoranthene-d10	19.08	212	10417	0.23	ng/ul	0.00
Target Compounds				Ovalue		
5) Naphthalene	10.48	128	1190	0.03	ng/ul#	77
7) 2-Methylnaphthalene	12.09	142	588	0.02	ng/ul	100
9) Acenaphthylene	14.01	152	2722	0.08	ng/ul#	94
11) Fluorene	15.34	166	1111	0.03	ng/ul#	76
14) Phenanthrene	17.07	178	25444m	0.43	ng/ul	
15) Anthracene	17.16	178	2970	0.05	ng/ul#	74
17) Fluoranthene	19.10	202	44161	0.68	ng/ul	95
19) Pyrene	19.46	202	40527	0.70	ng/ul	96
20) Benzo(a)anthracene	21.22	228	13239m	0.30	ng/ul	
21) Chrysene	21.28	228	16758	0.38	ng/ul	96
23) Benzo(b)fluoranthene	22.79	252	20701m	0.42	ng/ul	
24) Benzo(k)fluoranthene	22.81	252	8179m	0.18	ng/ul	
25) Benzo(a)pyrene	23.35	252	15285	0.34	ng/ul#	91
26) Indeno(1,2,3-cd)pyrene	25.66	276	10956	0.22	ng/ul#	87
27) Dibenzo(a,h)anthracene	25.68	278	2756	0.07	ng/ul#	35
28) Benzo(a,h,i)perylene	26.34	276	9865	0.23	ng/ul	97

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

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