

Data Path : Z:\HPCHEM1\BNA M\DATA\BM070716\
 Data File : BM006297.D
 Acq On : 06 Jul 2016 23:52
 Operator : UM/SJ
 Sample : H3811-11
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4TV9

Manual Integrations

umangi
 7/7/2016 6:05:34 PM

Quant Time: Jul 07 11:34:11 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	4991	0.40	ng/ul	0.00
4) Naphthalene-d8	10.43	136	19978	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.30	164	11657	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.04	188	25024m	0.40	ng/ul	0.00
18) Chrysene-d12	21.23	240	15990	0.40	ng/ul	0.00
22) Perylene-d12	23.45	264	11754m	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.17	96	9190	4.36	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.02	152	7704	0.33	ng/ul	0.00
16) Fluoranthene-d10	19.08	212	16232	0.31	ng/ul	0.00
Target Compounds				Ovalue		
5) Naphthalene	10.48	128	2706	0.06	ng/ul#	88
7) 2-Methylnaphthalene	12.09	142	2422	0.08	ng/ul	100
9) Acenaphthylene	14.01	152	5126	0.13	ng/ul#	93
11) Fluorene	15.34	166	1571	0.04	ng/ul#	73
14) Phenanthrene	17.07	178	41658m	0.60	ng/ul	
15) Anthracene	17.16	178	4279	0.07	ng/ul#	78
17) Fluoranthene	19.10	202	81066m	1.06	ng/ul	
19) Pyrene	19.48	202	72083	1.08	ng/ul#	90
20) Benzo(a)anthracene	21.21	228	20403	0.39	ng/ul#	89
21) Chrysene	21.27	228	32842	0.65	ng/ul#	90
23) Benzo(b)fluoranthene	22.79	252	32126m	0.76	ng/ul	
24) Benzo(k)fluoranthene	22.81	252	11370m	0.29	ng/ul	
25) Benzo(a)pyrene	23.35	252	20266m	0.52	ng/ul	
26) Indeno(1,2,3-cd)pyrene	25.66	276	16141m	0.37	ng/ul	
27) Dibenzo(a,h)anthracene	25.66	278	4260m	0.13	ng/ul	
28) Benzo(a,h,i)perylene	26.34	276	15322m	0.42	ng/ul	

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

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