

Data Path : Z:\HPCHEM1\BNA M\DATA\BM070716\
 Data File : BM006292.D
 Acq On : 06 Jul 2016 20:51
 Operator : UM/SJ
 Sample : H3811-06
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4TV6

Manual Integrations

umangi
 7/7/2016 6:05:09 PM

Quant Time: Jul 07 11:10:42 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	4497	0.40	ng/ul	0.00
4) Naphthalene-d8	10.43	136	17694	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.30	164	10248	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.04	188	22074m	0.40	ng/ul	0.00
18) Chrysene-d12	21.24	240	15079m	0.40	ng/ul	0.00
22) Perylene-d12	23.45	264	12733	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.17	96	7387	3.89	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.02	152	6102	0.29	ng/ul	0.00
16) Fluoranthene-d10	19.07	212	13337	0.29	ng/ul	-0.01
Target Compounds				Ovalue		
5) Naphthalene	10.48	128	1189	0.03	ng/ul#	77
7) 2-Methylnaphthalene	12.09	142	629	0.02	ng/ul	99
9) Acenaphthylene	14.01	152	2825	0.08	ng/ul#	94
10) Acenaphthene	14.35	153	962	0.03	ng/ul#	67
11) Fluorene	15.34	166	1292	0.03	ng/ul#	80
14) Phenanthrene	17.07	178	34429m	0.56	ng/ul	
15) Anthracene	17.16	178	5769	0.10	ng/ul#	85
17) Fluoranthene	19.11	202	70868	1.05	ng/ul	94
19) Pyrene	19.47	202	62727	0.99	ng/ul	96
20) Benzo(a)anthracene	21.22	228	24240m	0.50	ng/ul	
21) Chrysene	21.27	228	27197	0.57	ng/ul	96
23) Benzo(b)fluoranthene	22.78	252	34346m	0.75	ng/ul	
24) Benzo(k)fluoranthene	22.82	252	7894m	0.19	ng/ul	
25) Benzo(a)pyrene	23.35	252	21508	0.51	ng/ul	98
26) Indeno(1,2,3-cd)pyrene	25.66	276	15840	0.34	ng/ul#	81
27) Dibenzo(a,h)anthracene	25.67	278	3997	0.11	ng/ul#	15
28) Benzo(g,h,i)perylene	26.34	276	14129	0.35	ng/ul	96

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

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