

Data Path : Z:\HPCHEM1\BNA M\DATA\BM070716\
 Data File : BM006288.D
 Acq On : 06 Jul 2016 18:26
 Operator : UM/SJ
 Sample : H3811-02
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4TV2

Manual Integrations

umangi
 7/7/2016 6:04:27 PM

Quant Time: Jul 07 10:49:36 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	4482	0.40	ng/ul	0.00
4) Naphthalene-d8	10.43	136	17624	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.30	164	9949	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.04	188	19292m	0.40	ng/ul	0.00
18) Chrysene-d12	21.24	240	12710	0.40	ng/ul	0.00
22) Perylene-d12	23.45	264	10764m	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.17	96	8142	4.30	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.02	152	6814	0.33	ng/ul	0.00
16) Fluoranthene-d10	19.07	212	12029	0.30	ng/ul	-0.01
Target Compounds				Ovalue		
5) Naphthalene	10.48	128	1287	0.03	ng/ul#	79
7) 2-Methylnaphthalene	12.09	142	831	0.03	ng/ul	97
9) Acenaphthylene	14.01	152	1358	0.04	ng/ul#	90
14) Phenanthrene	17.07	178	15040m	0.28	ng/ul	
15) Anthracene	17.16	178	2771	0.06	ng/ul#	72
17) Fluoranthene	19.11	202	27969	0.48	ng/ul	94
19) Pyrene	19.47	202	25434	0.48	ng/ul	96
20) Benzo(a)anthracene	21.22	228	10181	0.25	ng/ul#	91
21) Chrysene	21.27	228	12513	0.31	ng/ul	94
23) Benzo(b)fluoranthene	22.78	252	16651m	0.43	ng/ul	
24) Benzo(k)fluoranthene	22.83	252	5201m	0.15	ng/ul	
25) Benzo(a)pyrene	23.35	252	9889m	0.28	ng/ul	
26) Indeno(1,2,3-cd)pyrene	25.66	276	7938m	0.20	ng/ul	
28) Benzo(a,h,i)perylene	26.35	276	7759m	0.23	ng/ul	

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

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