

Data Path : Z:\HPCHEM1\BNA M\DATA\BM070816\
 Data File : BM006342.D
 Acq On : 08 Jul 2016 05:32
 Operator : UM/SJ
 Sample : H3811-10DL 5X
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4TV8DL

Manual Integrations

umangi
 7/8/2016 7:14:30 PM

Quant Time: Jul 08 08:51:05 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM070716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 08 08:04:54 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	4718	0.40	ng/ul	0.00
4) Naphthalene-d8	10.43	136	18417	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.28	164	10502	0.40	ng/ul	-0.02
12) Phenanthrene-d10	17.04	188	23106	0.40	ng/ul	0.00
18) Chrysene-d12	21.24	240	15673	0.40	ng/ul	0.00
22) Perylene-d12	23.45	264	13941	0.40	ng/ul	0.01
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.17	96	1310	0.66	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.02	152	984	0.05	ng/ul	0.00
16) Fluoranthene-d10	19.08	212	3208	0.07	ng/ul	0.00
Target Compounds				Ovalue		
9) Acenaphthylene	14.01	152	1432	0.04	ng/ul#	92
11) Fluorene	15.34	166	816	0.02	ng/ul#	81
14) Phenanthrene	17.07	178	15041	0.24	ng/ul#	93
15) Anthracene	17.16	178	3622m	0.06	ng/ul	
17) Fluoranthene	19.10	202	29849	0.42	ng/ul	97
19) Pyrene	19.46	202	27377	0.42	ng/ul	98
20) Benzo(a)anthracene	21.22	228	13056	0.26	ng/ul#	90
21) Chrysene	21.28	228	12752	0.26	ng/ul	94
23) Benzo(b)fluoranthene	22.79	252	15800m	0.32	ng/ul	
24) Benzo(k)fluoranthene	22.81	252	6070m	0.13	ng/ul	
25) Benzo(a)pyrene	23.35	252	11497	0.25	ng/ul#	92
26) Indeno(1,2,3-cd)pyrene	25.66	276	7994m	0.16	ng/ul	
27) Dibenzo(a,h)anthracene	25.66	278	1942	0.05	ng/ul#	50
28) Benzo(g,h,i)perylene	26.34	276	7943	0.18	ng/ul	96

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

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