

Data Path : Z:\HPCHEM1\BNA M\DATA\BM070816\
 Data File : BM006338.D
 Acq On : 08 Jul 2016 03:07
 Operator : UM/SJ
 Sample : H3811-14DL 5X
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4TW2DL

Manual Integrations

umangi
 7/8/2016 7:14:21 PM

Quant Time: Jul 08 08:29:49 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	6737	0.40	ng/ul	0.00
4) Naphthalene-d8	10.43	136	26802	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.30	164	15186	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.04	188	36009	0.40	ng/ul	0.00
18) Chrysene-d12	21.23	240	28344	0.40	ng/ul	-0.01
22) Perylene-d12	23.44	264	25741	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.17	96	1623	0.57	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.02	152	1152	0.04	ng/ul	0.00
16) Fluoranthene-d10	19.08	212	3844	0.05	ng/ul	0.00
Target Compounds				Ovalue		
9) Acenaphthylene	14.01	152	9687	0.19	ng/ul#	93
11) Fluorene	15.34	166	2682	0.05	ng/ul#	81
14) Phenanthrene	17.07	178	48988	0.49	ng/ul#	92
15) Anthracene	17.16	178	7043m	0.08	ng/ul	
17) Fluoranthene	19.10	202	105324	0.96	ng/ul	98
19) Pyrene	19.46	202	102555	0.86	ng/ul	99
20) Benzo(a)anthracene	21.22	228	39773	0.43	ng/ul#	86
21) Chrysene	21.27	228	41991	0.47	ng/ul#	91
23) Benzo(b)fluoranthene	22.79	252	49197m	0.53	ng/ul	
24) Benzo(k)fluoranthene	22.81	252	16052m	0.19	ng/ul	
25) Benzo(a)pyrene	23.35	252	34476	0.40	ng/ul	98
26) Indeno(1,2,3-cd)pyrene	25.65	276	27026	0.28	ng/ul#	94
27) Dibenzo(a,h)anthracene	25.66	278	6032m	0.08	ng/ul	
28) Benzo(g,h,i)perylene	26.33	276	23451	0.29	ng/ul	95

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

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