

Data Path : Z:\HPCHEM1\BNA M\DATA\BM070816\
 Data File : BM006340.D
 Acq On : 08 Jul 2016 04:19
 Operator : UM/SJ
 Sample : H3811-16DL 5X
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4TW4DL

Manual Integrations

umangi
 7/8/2016 7:14:25 PM

Quant Time: Jul 08 08:39:21 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	5913	0.40	ng/ul	0.00
4) Naphthalene-d8	10.43	136	22575	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.28	164	12354	0.40	ng/ul	-0.02
12) Phenanthrene-d10	17.04	188	26711	0.40	ng/ul	0.00
18) Chrysene-d12	21.24	240	17731	0.40	ng/ul	0.00
22) Perylene-d12	23.45	264	16145	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.17	96	899	0.36	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.02	152	882	0.03	ng/ul	0.00
16) Fluoranthene-d10	19.07	212	4183	0.07	ng/ul	0.00
Target Compounds				Ovalue		
9) Acenaphthylene	14.01	152	7413	0.18	ng/ul#	92
11) Fluorene	15.34	166	1647	0.04	ng/ul#	83
14) Phenanthrene	17.07	178	30176	0.41	ng/ul#	92
15) Anthracene	17.16	178	6934m	0.10	ng/ul	
17) Fluoranthene	19.11	202	76610	0.94	ng/ul	94
19) Pyrene	19.47	202	84995	1.14	ng/ul	95
20) Benzo(a)anthracene	21.22	228	34159	0.60	ng/ul#	90
21) Chrysene	21.27	228	39657	0.70	ng/ul	96
23) Benzo(b)fluoranthene	22.78	252	47465m	0.82	ng/ul	
24) Benzo(k)fluoranthene	22.83	252	11093m	0.21	ng/ul	
25) Benzo(a)pyrene	23.35	252	31219	0.58	ng/ul	100
26) Indeno(1,2,3-cd)pyrene	25.66	276	23439	0.39	ng/ul#	93
27) Dibenzo(a,h)anthracene	25.67	278	5440	0.12	ng/ul#	65
28) Benzo(g,h,i)perylene	26.34	276	19783	0.39	ng/ul	95

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

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