

Data Path : Z:\HPCHEM1\BNA M\DATA\BM032516\
 Data File : BM004784.D
 Acq On : 25 Mar 2016 19:06
 Operator : UM/SJ
 Sample : H1909-08
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4T44

Manual Integrations
 APPROVED

Sohil
 4/6/2016 1:28:49 PM

Quant Time: Apr 05 19:17:29 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM032516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 05 18:29:16 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	1361	0.40	ng/ul	0.00
4) Naphthalene-d8	10.62	136	6412	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.47	164	4024	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.21	188	9065m	0.40	ng/ul	0.00
18) Chrysene-d12	21.40	240	8710	0.40	ng/ul	0.00
22) Perylene-d12	23.69	264	7204m	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.32	96	6815	11.11	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.21	152	3138	0.35	ng/ul	0.00
16) Fluoranthene-d10	19.24	212	8074	0.38	ng/ul	0.00
Target Compounds				Ovalue		
5) Naphthalene	10.67	128	1187	0.08	ng/ul#	82
9) Acenaphthylene	14.18	152	1073	0.06	ng/ul#	88
11) Fluorene	15.51	166	317	0.02	ng/ul#	74
14) Phenanthrene	17.24	178	4553m	0.18	ng/ul	
15) Anthracene	17.33	178	1327	0.06	ng/ul#	78
17) Fluoranthene	19.27	202	13146	0.45	ng/ul	96
19) Pyrene	19.63	202	12266	0.41	ng/ul	97
20) Benzo(a)anthracene	21.38	228	5928	0.22	ng/ul#	89
21) Chrysene	21.43	228	6190	0.24	ng/ul	94
23) Benzo(b)fluoranthene	22.99	252	8680m	0.36	ng/ul	
24) Benzo(k)fluoranthene	23.01	252	3700m	0.16	ng/ul	
25) Benzo(a)pyrene	23.59	252	7739m	0.35	ng/ul	
26) Indeno(1,2,3-cd)pyrene	26.01	276	3593m	0.17	ng/ul	
27) Dibenzo(a,h)anthracene	26.02	278	932m	0.06	ng/ul	
28) Benzo(a,h,i)perylene	26.71	276	3369m	0.20	ng/ul	

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

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