

Data Path : Z:\HPCHEM1\BNA M\DATA\BM092815\
 Data File : BM002735.D
 Acq On : 28 Sep 2015 16:02
 Operator : UM/IZ
 Sample : G3798-17
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9G71

Manual Integrations
 APPROVED

MMDadoda
 9/29/2015 3:42:37 PM

Quant Time: Sep 29 05:28:06 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-SIM-BM092415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 29 00:32:23 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.74	152	10531	0.40	ng/ul	0.00
4) Naphthalene-d8	11.60	136	41151	0.40	ng/ul	0.01
8) Acenaphthene-d10	15.33	164	22327	0.40	ng/ul	0.00
12) Phenanthrene-d10	18.03	188	45828	0.40	ng/ul	0.00
18) Chrysene-d12	22.13	240	44847	0.40	ng/ul	0.00
22) Perylene-d12	24.73	264	42161	0.40	ng/ul	0.01
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.80	96	31509	2.90	ng/uL	0.00
6) 2-Methylnaphthalene-d10	13.13	152	8193m	0.14	ng/ul	0.00
16) Fluoranthene-d10	20.01	212	14662	0.13	ng/ul	0.01
Target Compounds				Ovalue		
5) Naphthalene	11.65	128	16028	0.14	ng/ul#	86
7) 2-Methylnaphthalene	13.20	142	9303	0.13	ng/ul	94
9) Acenaphthylene	15.06	152	12338	0.11	ng/ul#	89
10) Acenaphthene	15.38	153	5415	0.06	ng/ul	91
11) Fluorene	16.36	166	3858m	0.04	ng/ul	
14) Phenanthrene	18.08	178	9896	0.07	ng/ul#	78
15) Anthracene	18.17	178	25328m	0.18	ng/ul	
17) Fluoranthene	20.03	202	73274	0.45	ng/ul	92
19) Pyrene	20.40	202	59706	0.34	ng/ul#	88
20) Benzo(a)anthracene	22.10	228	18285	0.12	ng/ul#	71
21) Chrysene	22.16	228	15317	0.11	ng/ul#	56
26) Indeno(1,2,3-cd)pyrene	27.46	276	9495	0.06	ng/ul	97
28) Benzo(g,h,i)perylene	28.33	276	10053	0.08	ng/ul#	27

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

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