

Data Path : Z:\HPCHEM1\BNA M\DATA\BM092815\
 Data File : BM002744.D
 Acq On : 28 Sep 2015 21:30
 Operator : UM/IZ
 Sample : G3798-15
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9G69

Manual Integrations
 APPROVED

MMDadoda
 9/29/2015 3:43:41 PM

Quant Time: Sep 29 07:51:35 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 02:52:19 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.75	152	8862	0.40	ng/ul	0.02
4) Naphthalene-d8	11.60	136	34468	0.40	ng/ul	0.00
8) Acenaphthene-d10	15.33	164	18375	0.40	ng/ul	0.00
12) Phenanthrene-d10	18.05	188	38689m	0.40	ng/ul	0.02
18) Chrysene-d12	22.13	240	37520	0.40	ng/ul	0.00
22) Perylene-d12	24.73	264	36978m	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.80	96	17055	1.86	ng/uL	0.00
6) 2-Methylnaphthalene-d10	13.14	152	5232m	0.10	ng/ul	0.01
16) Fluoranthene-d10	20.01	212	9175	0.10	ng/ul	0.00
Target Compounds				Ovalue		
5) Naphthalene	11.65	128	4919	0.05	ng/ul#	95
7) 2-Methylnaphthalene	13.20	142	4212	0.07	ng/ul	100
9) Acenaphthylene	15.06	152	19706	0.21	ng/ul	98
10) Acenaphthene	15.38	153	3135	0.05	ng/ul#	90
11) Fluorene	16.36	166	5198m	0.07	ng/ul	
14) Phenanthrene	18.08	178	26687	0.21	ng/ul	98
15) Anthracene	18.17	178	6220m	0.05	ng/ul	
17) Fluoranthene	20.03	202	58714m	0.43	ng/ul	
19) Pyrene	20.40	202	53105	0.36	ng/ul#	93
20) Benzo(a)anthracene	22.10	228	31589	0.25	ng/ul#	87
21) Chrysene	22.16	228	30086	0.25	ng/ul#	77
23) Benzo(b)fluoranthene	23.93	252	45931m	0.34	ng/ul	
24) Benzo(k)fluoranthene	23.97	252	15076m	0.12	ng/ul	
25) Benzo(a)pyrene	24.62	252	31783m	0.26	ng/ul	
26) Indeno(1,2,3-cd)pyrene	27.47	276	23846m	0.17	ng/ul	
28) Benzo(a,h,i)perylene	28.34	276	20619m	0.18	ng/ul	

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

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