

Data Path : Z:\HPCHEM1\BNA M\DATA\BM070816\
 Data File : BM006335.D
 Acq On : 08 Jul 2016 00:43
 Operator : UM/SJ
 Sample : H3811-15
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4TW3

Manual Integrations
 APPROVED

umangi
 7/8/2016 7:14:15 PM

Quant Time: Jul 08 07:52:09 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 03:28:13 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	5028	0.40	ng/ul	0.00
4) Naphthalene-d8	10.43	136	20035	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.30	164	11653	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.04	188	23605m	0.40	ng/ul	0.00
18) Chrysene-d12	21.24	240	14775m	0.40	ng/ul	0.00
22) Perylene-d12	23.45	264	14565	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.17	96	9855	4.64	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.02	152	7764	0.33	ng/ul	0.00
16) Fluoranthene-d10	19.07	212	15752	0.32	ng/ul	-0.01
Target Compounds				Ovalue		
5) Naphthalene	10.46	128	11818	0.25	ng/ul	97
7) 2-Methylnaphthalene	12.09	142	6284	0.20	ng/ul	100
9) Acenaphthylene	14.01	152	60774	1.55	ng/ul	92
10) Acenaphthene	14.35	153	5893	0.15	ng/ul	89
11) Fluorene	15.34	166	24257m	0.55	ng/ul	
14) Phenanthrene	17.07	178	388843m	5.96	ng/ul	
15) Anthracene	17.16	178	84242m	1.39	ng/ul	
17) Fluoranthene	19.11	202	603410m	8.38	ng/ul	
19) Pyrene	19.47	202	799339	12.92	ng/ul	97
20) Benzo(a)anthracene	21.22	228	362076m	7.57	ng/ul	
21) Chrysene	21.27	228	363462	7.75	ng/ul	97
23) Benzo(b)fluoranthene	22.78	252	318074m	6.10	ng/ul	
24) Benzo(k)fluoranthene	22.82	252	77226m	1.59	ng/ul	
25) Benzo(a)pyrene	23.35	252	260296m	5.36	ng/ul	
26) Indeno(1,2,3-cd)pyrene	25.66	276	148219	2.75	ng/ul	97
27) Dibenzo(a,h)anthracene	25.66	278	44731	1.06	ng/ul	97
28) Benzo(g,h,i)perylene	26.34	276	136765	3.00	ng/ul#	94

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

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