

Data Path : Z:\HPCHEM1\BNA M\DATA\BM070816\
 Data File : BM006339.D
 Acq On : 08 Jul 2016 03:43
 Operator : UM/SJ
 Sample : H3811-15DL 5X
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4TW3DL

Manual Integrations

umangi
 7/8/2016 7:14:23 PM

Quant Time: Jul 08 08:35:21 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	4722	0.40	ng/ul	0.00
4) Naphthalene-d8	10.43	136	18840	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.28	164	11098	0.40	ng/ul	-0.02
12) Phenanthrene-d10	17.04	188	25977	0.40	ng/ul	0.00
18) Chrysene-d12	21.23	240	17984	0.40	ng/ul	-0.01
22) Perylene-d12	23.44	264	15080	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.17	96	1978	0.99	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.02	152	1488	0.07	ng/ul	0.00
16) Fluoranthene-d10	19.08	212	5399	0.10	ng/ul	0.00
Target Compounds				Ovalue		
5) Naphthalene	10.46	128	2330	0.05	ng/ul#	86
7) 2-Methylnaphthalene	12.09	142	1135	0.04	ng/ul	99
9) Acenaphthylene	14.01	152	14161	0.38	ng/ul#	92
10) Acenaphthene	14.35	153	1354	0.04	ng/ul#	89
11) Fluorene	15.34	166	5289m	0.13	ng/ul	
14) Phenanthrene	17.07	178	93853	1.31	ng/ul#	93
15) Anthracene	17.16	178	20899m	0.31	ng/ul	
17) Fluoranthene	19.10	202	176739	2.23	ng/ul	98
19) Pyrene	19.46	202	223961	2.97	ng/ul	99
20) Benzo(a)anthracene	21.22	228	98539	1.69	ng/ul#	89
21) Chrysene	21.27	228	94057	1.65	ng/ul#	89
23) Benzo(b)fluoranthene	22.79	252	69002m	1.28	ng/ul	
24) Benzo(k)fluoranthene	22.81	252	25498m	0.51	ng/ul	
25) Benzo(a)pyrene	23.35	252	59454m	1.18	ng/ul	
26) Indeno(1,2,3-cd)pyrene	25.65	276	31406	0.56	ng/ul#	95
27) Dibenzo(a,h)anthracene	25.66	278	9554m	0.22	ng/ul	
28) Benzo(g,h,i)perylene	26.33	276	28793	0.61	ng/ul	96

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

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