

Data Path : Z:\HPCHEM1\BNA_M\Data\BM061416\
 Data File : BM005782.D
 Acq On : 14 Jun 2016 12:49
 Operator : UM/SJ
 Sample : H3507-01DL 5X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JHGK1DL

Manual Integrations
 APPROVED

umangi
 6/14/2016 5:05:19 PM

Quant Time: Jun 14 16:05:16 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM061316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 13 19:45:38 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	6099	0.40	ng/ul	-0.02
4) Naphthalene-d8	10.54	136	26637	0.40	ng/ul	-0.01
8) Acenaphthene-d10	14.38	164	18782	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.12	188	35754	0.40	ng/ul	-0.02
18) Chrysene-d12	21.33	240	22962	0.40	ng/ul	0.00
22) Perylene-d12	23.58	264	26746	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.20	96	4524	1.52	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.14	152	1310	0.04	ng/ul	0.00
16) Fluoranthene-d10	19.16	212	3033	0.04	ng/ul	0.00
Target Compounds						
9) Acenaphthylene	14.11	152	3113	0.04	ng/ul#	93
11) Fluorene	15.44	166	3324	0.06	ng/ul#	93
14) Phenanthrene	17.18	178	34399	0.40	ng/ul	98
15) Anthracene	17.26	178	8439m	0.08	ng/ul	
17) Fluoranthene	19.20	202	49189	0.46	ng/ul	94
19) Pyrene	19.56	202	46732	0.59	ng/ul#	92
20) Benzo(a)anthracene	21.31	228	15754	0.29	ng/ul#	85
21) Chrysene	21.36	228	15895	0.19	ng/ul	92
23) Benzo(b)fluoranthene	22.90	252	16893m	0.23	ng/ul	
24) Benzo(k)fluoranthene	22.92	252	8456m	0.09	ng/ul	
25) Benzo(a)pyrene	23.48	252	11803	0.14	ng/ul#	83
26) Indeno(1,2,3-cd)pyrene	25.86	276	7111	0.07	ng/ul	94
28) Benzo(g,h,i)perylene	26.56	276	6527	0.08	ng/ul#	83

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

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