

Data Path : Z:\HPCHEM1\BNA M\DATA\BM121216\
 Data File : BM008265.D
 Acq On : 12 Dec 2016 16:08
 Operator : UM/SJ
 Sample : H5905-09DL 5X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA809DL

Manual Integrations
 APPROVED

umangi
 12/13/2016 6:12:38 PM

Quant Time: Dec 13 03:02:40 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 13 02:27:35 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	630	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	2102	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.33	164	1196	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	2081	0.40	ng/ul	-0.02
16) Chrysene-d12	21.27	240	2052	0.40	ng/ul	-0.01
20) Perylene-d12	23.51	264	1945	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.07	152	131	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.11	212	279m	0.05	ng/ul	-0.01
Target Compounds				Ovalue		
3) Naphthalene	10.51	128	1441	0.25	ng/ul#	92
5) 2-Methylnaphthalene	12.14	142	1589	0.42	ng/ul	94
7) Acenaphthylene	14.04	152	1209	0.22	ng/ul#	89
8) Acenaphthene	14.38	153	114	0.03	ng/ul#	83
9) Fluorene	15.38	166	104	0.02	ng/ul#	47
12) Phenanthrene	17.11	178	2510	0.39	ng/ul	97
13) Anthracene	17.21	178	1377m	0.23	ng/ul	
15) Fluoranthene	19.14	202	4216m	0.50	ng/ul	
17) Pyrene	19.51	202	5028	0.71	ng/ul	97
18) Benzo(a)anthracene	21.25	228	4906	0.78	ng/ul	95
19) Chrysene	21.30	228	7567	0.97	ng/ul	93
21) Benzo(b)fluoranthene	22.84	252	14545m	1.87	ng/ul	
22) Benzo(k)fluoranthene	22.88	252	3631m	0.45	ng/ul	
23) Benzo(a)pyrene	23.41	252	7255m	1.00	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.74	276	6823	0.76	ng/ul#	91
25) Dibenzo(a,h)anthracene	25.74	278	1822	0.25	ng/ul#	87
26) Benzo(a,h,i)perylene	26.44	276	5913	0.74	ng/ul#	91

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

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