

Data Path : Z:\HPCHEM1\BNA M\DATA\BM121216\
 Data File : BM008267.D
 Acq On : 12 Dec 2016 17:21
 Operator : UM/SJ
 Sample : H5905-04DL 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA806DL

Manual Integrations
 APPROVED

umangi
 12/13/2016 6:12:32 PM

Quant Time: Dec 13 03:07:51 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	541	0.40	ng/ul	0.00
2) Naphthalene-d8	10.48	136	1817	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.33	164	1034	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	1744	0.40	ng/ul	-0.02
16) Chrysene-d12	21.27	240	1526	0.40	ng/ul	-0.01
20) Perylene-d12	23.51	264	1499	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.07	152	102	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.11	212	224m	0.05	ng/ul	-0.01
Target Compounds						
					Ovalue	
3) Naphthalene	10.52	128	836	0.16	ng/ul#	82
5) 2-Methylnaphthalene	12.14	142	686	0.21	ng/ul	98
7) Acenaphthylene	14.04	152	594	0.13	ng/ul#	77
8) Acenaphthene	14.38	153	276	0.08	ng/ul#	92
9) Fluorene	15.38	166	195m	0.05	ng/ul	
12) Phenanthrene	17.11	178	2658	0.49	ng/ul	97
13) Anthracene	17.21	178	1031m	0.20	ng/ul	
15) Fluoranthene	19.14	202	5532	0.79	ng/ul	96
17) Pyrene	19.51	202	5584	1.05	ng/ul	98
18) Benzo(a)anthracene	21.25	228	3483m	0.74	ng/ul	
19) Chrysene	21.30	228	4235m	0.73	ng/ul	
21) Benzo(b)fluoranthene	22.84	252	7659m	1.28	ng/ul	
22) Benzo(k)fluoranthene	22.88	252	2000m	0.32	ng/ul	
23) Benzo(a)pyrene	23.41	252	3677	0.66	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.76	276	3047	0.44	ng/ul#	90
25) Dibenzo(a,h)anthracene	25.76	278	839	0.15	ng/ul#	69
26) Benzo(a,h,i)perylene	26.44	276	2398	0.39	ng/ul	94

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

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