

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112316\
 Data File : BM008063.D
 Acq On : 24 Nov 2016 08:32
 Operator : UM/SJ
 Sample : H5701-24
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4WG2

Manual Integrations
 APPROVED

Umangi
 11/28/2016 5:29:04 PM

Quant Time: Nov 25 01:22:22 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 24 04:11:15 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	844	0.40	ng/ul	-0.03
2) Naphthalene-d8	10.49	136	2936	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.35	164	1810	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	3647m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.29	240	3348	0.40	ng/ul	-0.01
20) Perylene-d12	23.52	264	2881m	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.09	152	1151	0.28	ng/ul	-0.03
14) Fluoranthene-d10	19.13	212	2631	0.25	ng/ul	-0.01
Target Compounds					Ovalue	
3) Naphthalene	10.54	128	237	0.03	ng/ul#	50
7) Acenaphthylene	14.06	152	649	0.08	ng/ul#	84
8) Acenaphthene	14.40	153	275	0.04	ng/ul#	94
9) Fluorene	15.39	166	448	0.06	ng/ul#	81
12) Phenanthrene	17.12	178	6248	0.55	ng/ul	97
13) Anthracene	17.23	178	2029	0.19	ng/ul#	93
15) Fluoranthene	19.16	202	13925	0.95	ng/ul	94
17) Pyrene	19.52	202	12455	1.07	ng/ul	98
18) Benzo(a)anthracene	21.27	228	6719	0.65	ng/ul	97
19) Chrysene	21.32	228	6638	0.52	ng/ul	97
21) Benzo(b)fluoranthene	22.86	252	8737m	0.76	ng/ul	
22) Benzo(k)fluoranthene	22.89	252	3475m	0.29	ng/ul	
23) Benzo(a)pyrene	23.42	252	6986m	0.65	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.78	276	4080m	0.31	ng/ul	
25) Dibenzo(a,h)anthracene	25.79	278	1180m	0.11	ng/ul	
26) Benzo(a,h,i)perylene	26.46	276	3398m	0.29	ng/ul	

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

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