

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030817\
 Data File : BM009131.D
 Acq On : 09 Mar 2017 06:22
 Operator : SJ/MA
 Sample : I2069-06
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AG7

Manual Integrations
 APPROVED

mohammad
 3/9/2017 2:02:46 PM

Quant Time: Mar 09 07:06:52 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM030317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 09 02:44:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	294	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	929	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.52	164	566	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	1269m	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	1390	0.40	ng/ul	-0.01
20) Perylene-d12	23.78	264	1309m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	379	0.25	ng/ul	0.00
14) Fluoranthene-d10	19.30	212	1045	0.26	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.73	128	323	0.12	ng/ul#	64
5) 2-Methylnaphthalene	12.34	142	174	0.09	ng/ul	100
8) Acenaphthene	14.59	153	427	0.19	ng/ul	93
9) Fluorene	15.58	166	681m	0.26	ng/ul	
12) Phenanthrene	17.31	178	9289m	2.33	ng/ul	
13) Anthracene	17.40	178	2615	0.66	ng/ul	98
15) Fluoranthene	19.33	202	12664	2.44	ng/ul	93
17) Pyrene	19.69	202	10768	2.25	ng/ul	97
18) Benzo(a)anthracene	21.42	228	5900	1.28	ng/ul	95
19) Chrysene	21.48	228	4519	1.03	ng/ul	96
21) Benzo(b)fluoranthene	23.07	252	6373m	1.24	ng/ul	
22) Benzo(k)fluoranthene	23.11	252	2206m	0.49	ng/ul	
23) Benzo(a)pyrene	23.67	252	4632m	1.00	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.15	276	2723m	0.46	ng/ul	
25) Dibenzo(a,h)anthracene	26.16	278	797m	0.16	ng/ul	
26) Benzo(g,h,i)perylene	26.89	276	2299m	0.46	ng/ul	

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

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