

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031117\
 Data File : BM009214.D
 Acq On : 11 Mar 2017 18:53
 Operator : SJ/MA
 Sample : I2072-08DL 5X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AK3DL

Manual Integrations
 APPROVED

mohammad
 3/13/2017 5:14:07 PM

Quant Time: Mar 13 12:20:52 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM030317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 13 11:21:07 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	212	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	659	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	362	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	818	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	928	0.40	ng/ul	0.00
20) Perylene-d12	23.77	264	964	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	44	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	116	0.05	ng/ul	0.00
Target Compounds						
3) Naphthalene	10.73	128	499	0.26	ng/ul#	78
5) 2-Methylnaphthalene	12.34	142	388	0.28	ng/ul	98
12) Phenanthrene	17.30	178	528	0.21	ng/ul#	79
15) Fluoranthene	19.31	202	906	0.27	ng/ul	91
17) Pyrene	19.68	202	1034	0.32	ng/ul#	87
18) Benzo(a)anthracene	21.42	228	720	0.23	ng/ul#	89
19) Chrysene	21.47	228	752m	0.26	ng/ul	
21) Benzo(b)fluoranthene	23.07	252	1077m	0.28	ng/ul	
22) Benzo(k)fluoranthene	23.09	252	461m	0.14	ng/ul	
23) Benzo(a)pyrene	23.67	252	687	0.20	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	26.14	276	488	0.11	ng/ul#	86
26) Benzo(g,h,i)perylene	26.88	276	443	0.12	ng/ul#	70

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

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