

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080318\
 Data File : BE097183.D
 Acq On : 3 Aug 2018 17:18
 Operator : SJ/JU
 Sample : J4171-05
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Aug 03 23:54:06 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE080118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 03 23:39:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1510	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	7960	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	5605	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	12629m	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	13081m	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	11635m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	5844	0.46	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	17222m	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.19	128	4716	0.245	ng/ul#	83
5) 2-Methylnaphthalene	11.80	142	2647	0.208	ng/ul	94
8) Acenaphthene	14.09	153	1118	0.064	ng/ul#	87
9) Fluorene	15.08	166	490	0.024	ng/ul#	51
12) Phenanthrene	16.81	178	7167m	0.223	ng/ul	
13) Anthracene	16.90	178	1150m	0.038	ng/ul	
15) Fluoranthene	18.84	202	14896m	0.350	ng/ul	
17) Pyrene	19.21	202	18744m	0.535	ng/ul	
18) Benzo(a)anthracene	20.96	228	6715m	0.180	ng/ul	
19) Chrysene	21.02	228	9389m	0.244	ng/ul	
21) Benzo(b)fluoranthene	22.54	252	10165m	0.321	ng/ul	
22) Benzo(k)fluoranthene	22.56	252	6872m	0.201	ng/ul	
23) Benzo(a)pyrene	23.12	252	8028	0.253	ng/ul#	74
24) Indeno(1,2,3-cd)pyrene	25.50	276	5971	0.150	ng/ul#	78
26) Benzo(a,h,i)perylene	26.21	276	6418	0.187	ng/ul#	87

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

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