

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080318\
 Data File : BE097186.D
 Acq On : 3 Aug 2018 19:17
 Operator : SJ/JU
 Sample : J4171-20
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Aug 04 00:08:56 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	1524	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	7842	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	5237	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	11818m	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	12339m	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	12336m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	4723	0.37	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	13530m	0.33	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	1099	0.058	ng/ul#	89
5) 2-Methylnaphthalene	11.80	142	503	0.040	ng/ul	99
7) Acenaphthylene	13.74	152	972	0.043	ng/ul#	72
8) Acenaphthene	14.08	153	2527	0.154	ng/ul	96
9) Fluorene	15.08	166	3149	0.167	ng/ul	91
12) Phenanthrene	16.81	178	143746m	4.771	ng/ul	
13) Anthracene	16.90	178	24646m	0.877	ng/ul	
15) Fluoranthene	18.84	202	386551m	9.707	ng/ul	
17) Pyrene	19.21	202	411281	12.439	ng/ul#	80
18) Benzo(a)anthracene	20.96	228	176772m	5.014	ng/ul	
19) Chrysene	21.02	228	198163	5.454	ng/ul	99
21) Benzo(b)fluoranthene	22.54	252	246705m	7.338	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	91830m	2.533	ng/ul	
23) Benzo(a)pyrene	23.12	252	194472m	5.780	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.51	276	143779m	3.415	ng/ul	
25) Dibenzo(a,h)anthracene	25.52	278	32913m	0.955	ng/ul	
26) Benzo(a,h,i)perylene	26.22	276	135717m	3.730	ng/ul	

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

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