

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080318\
 Data File : BE097190.D
 Acq On : 3 Aug 2018 21:47
 Operator : SJ/JU
 Sample : J4171-21
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Aug 04 00:29:42 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	7.40	152	1392	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	7447	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	4892	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	11132m	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	11952m	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	11198m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	4387	0.37	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	13453m	0.35	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	795	0.044	ng/ul#	83
5) 2-Methylnaphthalene	11.81	142	338	0.028	ng/ul	98
7) Acenaphthylene	13.74	152	2698	0.128	ng/ul	97
9) Fluorene	15.08	166	956	0.054	ng/ul#	90
12) Phenanthrene	16.81	178	37217m	1.311	ng/ul	
13) Anthracene	16.90	178	1656m	0.063	ng/ul	
15) Fluoranthene	18.84	202	53108m	1.416	ng/ul	
17) Pyrene	19.21	202	54572	1.704	ng/ul#	79
18) Benzo(a)anthracene	20.96	228	14366m	0.421	ng/ul	
19) Chrysene	21.01	228	24686	0.701	ng/ul	96
21) Benzo(b)fluoranthene	22.54	252	28558m	0.936	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	11000m	0.334	ng/ul	
23) Benzo(a)pyrene	23.12	252	20548m	0.673	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.51	276	16604	0.434	ng/ul#	72
25) Dibenzo(a,h)anthracene	25.53	278	3532	0.113	ng/ul#	62
26) Benzo(a,h,i)perylene	26.21	276	17052	0.516	ng/ul	94

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

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