

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080318\
 Data File : BE097187.D
 Acq On : 3 Aug 2018 19:53
 Operator : SJ/JU
 Sample : J4171-18
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Aug 04 00:14:34 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	1405	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	7388	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	4870	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	11351m	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	11722m	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	11347m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	3305	0.28	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	9411m	0.24	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	700	0.039	ng/ul#	79
5) 2-Methylnaphthalene	11.80	142	780	0.066	ng/ul	97
12) Phenanthrene	16.81	178	12619m	0.436	ng/ul	
13) Anthracene	16.90	178	2376m	0.088	ng/ul	
15) Fluoranthene	18.84	202	37231m	0.973	ng/ul	
17) Pyrene	19.21	202	44511	1.417	ng/ul#	81
18) Benzo(a)anthracene	20.96	228	20855m	0.623	ng/ul	
19) Chrysene	21.01	228	27194m	0.788	ng/ul	
21) Benzo(b)fluoranthene	22.55	252	39056m	1.263	ng/ul	
22) Benzo(k)fluoranthene	22.57	252	17418m	0.522	ng/ul	
23) Benzo(a)pyrene	23.12	252	28091m	0.908	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.52	276	34438m	0.889	ng/ul	
25) Dibenzo(a,h)anthracene	25.53	278	6717m	0.212	ng/ul	
26) Benzo(g,h,i)perylene	26.22	276	37751m	1.128	ng/ul	

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

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