

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080318\
 Data File : BE097185.D
 Acq On : 3 Aug 2018 18:38
 Operator : SJ/JU
 Sample : J4171-16
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Aug 04 00:04:10 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	1479	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	7720	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	4844	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	11731m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	12488m	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	12410m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	3350	0.27	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	9823m	0.24	ng/ul	0.00
Target Compounds				Ovalue		
8) Acenaphthene	14.08	153	554	0.037	ng/ul#	93
9) Fluorene	15.08	166	834	0.048	ng/ul#	90
12) Phenanthrene	16.81	178	43541m	1.456	ng/ul	
13) Anthracene	16.90	178	6568m	0.235	ng/ul	
15) Fluoranthene	18.84	202	139645m	3.533	ng/ul	
17) Pyrene	19.21	202	150321	4.492	ng/ul#	79
18) Benzo(a)anthracene	20.96	228	64828m	1.817	ng/ul	
19) Chrysene	21.01	228	76727	2.087	ng/ul	98
21) Benzo(b)fluoranthene	22.54	252	101952m	3.014	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	30820m	0.845	ng/ul	
23) Benzo(a)pyrene	23.11	252	77705m	2.296	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.50	276	55870m	1.319	ng/ul	
25) Dibenzo(a,h)anthracene	25.52	278	12801m	0.369	ng/ul	
26) Benzo(g,h,i)perylene	26.21	276	53200m	1.453	ng/ul	

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

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