

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080418\
 Data File : BE097206.D
 Acq On : 4 Aug 2018 17:58
 Operator : SJ/JU
 Sample : J4206-04
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Aug 06 02:15:10 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 : Mon Aug 06 00:53:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1333	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	7059	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	4498	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	10572m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	10670m	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	9945m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	2995	0.26	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	8721m	0.24	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	728	0.043	ng/ul#	82
5) 2-Methylnaphthalene	11.80	142	466	0.041	ng/ul	96
7) Acenaphthylene	13.73	152	690	0.035	ng/ul#	78
8) Acenaphthene	14.08	153	295	0.021	ng/ul#	90
9) Fluorene	15.08	166	659	0.041	ng/ul#	87
12) Phenanthrene	16.81	178	9108m	0.338	ng/ul	
13) Anthracene	16.90	178	1487m	0.059	ng/ul	
15) Fluoranthene	18.84	202	11933m	0.335	ng/ul	
17) Pyrene	19.20	202	14434	0.505	ng/ul#	82
18) Benzo(a)anthracene	20.96	228	5152	0.169	ng/ul#	85
19) Chrysene	21.01	228	6243	0.199	ng/ul	94
21) Benzo(b)fluoranthene	22.54	252	6840m	0.252	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	2007m	0.069	ng/ul	
23) Benzo(a)pyrene	23.11	252	5507	0.203	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.50	276	3479	0.102	ng/ul#	75
25) Dibenzo(a,h)anthracene	25.52	278	814	0.029	ng/ul#	45
26) Benzo(a,h,i)perylene	26.21	276	3680	0.125	ng/ul	93

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

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