

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080418\
 Data File : BE097212.D
 Acq On : 4 Aug 2018 22:21
 Operator : SJ/JU
 Sample : J4206-06
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Aug 06 04:20:10 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 03:46:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1513	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	8150	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	4987	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	11377m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	11882m	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	10801m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	3059	0.23	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	8810m	0.23	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	1114	0.056	ng/ul#	90
5) 2-Methylnaphthalene	11.80	142	501	0.038	ng/ul	96
7) Acenaphthylene	13.73	152	1316	0.061	ng/ul#	87
8) Acenaphthene	14.08	153	5506	0.353	ng/ul	95
9) Fluorene	15.08	166	6012	0.335	ng/ul	95
12) Phenanthrene	16.81	178	205376m	7.081	ng/ul	
13) Anthracene	16.90	178	30237m	1.118	ng/ul	
15) Fluoranthene	18.84	202	456377m	11.905	ng/ul	
17) Pyrene	19.21	202	482284	15.148	ng/ul#	79
18) Benzo(a)anthracene	20.96	228	193722m	5.706	ng/ul	
19) Chrysene	21.01	228	229102	6.548	ng/ul	97
21) Benzo(b)fluoranthene	22.54	252	283565m	9.632	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	105660m	3.329	ng/ul	
23) Benzo(a)pyrene	23.12	252	222878m	7.566	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.51	276	163945m	4.447	ng/ul	
25) Dibenzo(a,h)anthracene	25.52	278	38068m	1.262	ng/ul	
26) Benzo(a,h,i)perylene	26.22	276	163207m	5.123	ng/ul	

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

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