

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080418\
 Data File : BE097207.D
 Acq On : 4 Aug 2018 18:37
 Operator : SJ/JU
 Sample : J4206-02
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Aug 06 03:10:00 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	1429	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	7656	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	4859	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	11270m	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	11846m	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	10587m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	2284	0.19	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	7608m	0.20	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	3447	0.186	ng/ul#	95
5) 2-Methylnaphthalene	11.80	142	689	0.056	ng/ul	98
7) Acenaphthylene	13.73	152	1733	0.082	ng/ul#	81
8) Acenaphthene	14.08	153	584	0.038	ng/ul#	88
9) Fluorene	15.08	166	790	0.045	ng/ul#	83
12) Phenanthrene	16.81	178	17523m	0.610	ng/ul	
13) Anthracene	16.90	178	1741m	0.065	ng/ul	
15) Fluoranthene	18.84	202	23989m	0.632	ng/ul	
17) Pyrene	19.21	202	31006	0.977	ng/ul#	81
18) Benzo(a)anthracene	20.96	228	9703m	0.287	ng/ul	
19) Chrysene	21.01	228	16565m	0.475	ng/ul	
21) Benzo(b)fluoranthene	22.54	252	18172m	0.630	ng/ul	
22) Benzo(k)fluoranthene	22.57	252	6375m	0.205	ng/ul	
23) Benzo(a)pyrene	23.11	252	14638m	0.507	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.50	276	10260m	0.284	ng/ul	
25) Dibenzo(a,h)anthracene	25.52	278	2408m	0.081	ng/ul	
26) Benzo(a,h,i)perylene	26.21	276	10951m	0.351	ng/ul	

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

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