

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080418\
 Data File : BE097204.D
 Acq On : 4 Aug 2018 16:44
 Operator : SJ/JU
 Sample : J4206-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Aug 06 02:00:58 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	1290	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	7025	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	4411	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	10296m	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	10457m	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	9689m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	3156	0.28	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	9985m	0.28	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	507	0.030	ng/ul#	76
5) 2-Methylnaphthalene	11.80	142	459	0.041	ng/ul	98
12) Phenanthrene	16.81	178	1192m	0.045	ng/ul	
15) Fluoranthene	18.84	202	1589m	0.046	ng/ul	
17) Pyrene	19.21	202	2118	0.076	ng/ul	94
18) Benzo(a)anthracene	20.96	228	938m	0.031	ng/ul	
19) Chrysene	21.01	228	1293m	0.042	ng/ul	
21) Benzo(b)fluoranthene	22.54	252	1483m	0.056	ng/ul	
22) Benzo(k)fluoranthene	22.57	252	693m	0.024	ng/ul	
23) Benzo(a)pyrene	23.11	252	989m	0.037	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.50	276	815	0.025	ng/ul#	89
26) Benzo(g,h,i)perylene	26.20	276	929	0.033	ng/ul#	71

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

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