

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080418\
 Data File : BE097203.D
 Acq On : 4 Aug 2018 16:05
 Operator : SJ/JU
 Sample : J4265-03
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Aug 06 01:50:15 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	1311	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	7115	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	4495	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	10398m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	10040m	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	9437m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	3357	0.29	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	9935m	0.28	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	16.81	178	970m	0.037	ng/ul	
13) Anthracene	16.90	178	945m	0.038	ng/ul	
15) Fluoranthene	18.84	202	1571m	0.045	ng/ul	
17) Pyrene	19.21	202	1657	0.062	ng/ul#	90
18) Benzo(a)anthracene	20.96	228	1550	0.054	ng/ul#	87
19) Chrysene	21.01	228	1644	0.056	ng/ul	93
21) Benzo(b)fluoranthene	22.54	252	1453	0.056	ng/ul	93
22) Benzo(k)fluoranthene	22.58	252	1527	0.055	ng/ul#	75
23) Benzo(a)pyrene	23.11	252	1338	0.052	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.51	276	1823	0.057	ng/ul#	82
25) Dibenzo(a,h)anthracene	25.53	278	1650	0.063	ng/ul#	81
26) Benzo(g,h,i)perylene	26.21	276	1564	0.056	ng/ul#	88

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

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