

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080418\
 Data File : BE097205.D
 Acq On : 4 Aug 2018 17:22
 Operator : SJ/JU
 Sample : J4206-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 JJ1J7

Manual Integrations
 APPROVED

Sohil
 8/6/2018 12:50:53 PM

Quant Time: Aug 14 00:51:18 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	1326	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	7212	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	4454	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	10538m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	10784m	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	9966m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	2552	0.22	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	7740m	0.21	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	16.81	178	2090m	0.078	ng/ul	
15) Fluoranthene	18.84	202	2342m	0.066	ng/ul	
17) Pyrene	19.21	202	2682m	0.093	ng/ul	
18) Benzo(a)anthracene	20.96	228	1022m	0.033	ng/ul	
19) Chrysene	21.01	228	1468m	0.046	ng/ul	
21) Benzo(b)fluoranthene	22.53	252	1835m	0.068	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	1000m	0.034	ng/ul	
23) Benzo(a)pyrene	23.11	252	1283	0.047	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	25.50	276	1435	0.042	ng/ul#	86
26) Benzo(g,h,i)perylene	26.20	276	1533	0.052	ng/ul#	85

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

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