

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110118\
 Data File : BN003390.D
 Acq On : 01 Nov 2018 21:56
 Operator : MJ/SJ
 Sample : J5701-12
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A40F9

Manual Integrations
 APPROVED

Sohil
 11/2/2018 4:14:14 PM

Quant Time: Nov 13 13:25:56 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN110118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 02 04:49:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	13456	0.40	ng/ul	0.00
2) Naphthalene-d8	10.66	136	53384	0.40	ng/ul	0.01
6) Acenaphthene-d10	14.49	164	30679	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.24	188	436839	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	59273	0.40	ng/ul	0.00
20) Perylene-d12	23.73	264	53289	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	18511	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	41897	0.03	ng/ul	0.06
Target Compounds				Ovalue		
9) Fluorene	15.54	166	111005	0.791	ng/ul#	92
15) Fluoranthene	19.35	202	87355	0.054	ng/ul	96
17) Pyrene	19.68	202	72969	0.338	ng/ul#	88
18) Benzo(a)anthracene	21.40	228	25674	0.138	ng/ul#	90
19) Chrysene	21.45	228	48299	0.229	ng/ul#	94
21) Benzo(b)fluoranthene	23.02	252	56785m	0.247	ng/ul	
22) Benzo(k)fluoranthene	23.07	252	19336m	0.087	ng/ul	
23) Benzo(a)pyrene	23.62	252	21884	0.114	ng/ul#	48
24) Indeno(1,2,3-cd)pyrene	26.08	276	18447	0.090	ng/ul#	82
25) Dibenzo(a,h)anthracene	26.09	278	4677	0.028	ng/ul#	1
26) Benzo(g,h,i)perylene	26.81	276	15447	0.088	ng/ul#	82

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

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