

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101118\
 Data File : BN003095.D
 Acq On : 11 Oct 2018 09:49
 Operator : MJ/SJ
 Sample : J5234-19DL 5X
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AK4DL

Manual Integrations
 APPROVED

Sohil
 10/11/2018 5:45:16 PM

Quant Time: Oct 11 13:41:08 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN100918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 11 08:10:30 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1384	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.40	136	5851	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.26	164	3553	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	8015	0.40	ng/ul	-0.02
16) Chrysene-d12	21.23	240	5674	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	4938	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	678	0.07	ng/ul	0.02
14) Fluoranthene-d10	19.06	212	1666	0.07	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.98	152	679	0.042	ng/ul#	78
9) Fluorene	15.32	166	747	0.053	ng/ul#	96
12) Phenanthrene	17.06	178	9750	0.436	ng/ul	100
13) Anthracene	17.15	178	1286	0.061	ng/ul#	85
15) Fluoranthene	19.09	202	16093	0.614	ng/ul	99
17) Pyrene	19.45	202	14114	0.640	ng/ul	97
18) Benzo(a)anthracene	21.22	228	5036m	0.276	ng/ul	
19) Chrysene	21.27	228	5259	0.284	ng/ul	96
21) Benzo(b)fluoranthene	22.79	252	5212m	0.293	ng/ul	
22) Benzo(k)fluoranthene	22.83	252	2155m	0.117	ng/ul	
23) Benzo(a)pyrene	23.34	252	3820	0.240	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.68	276	2471	0.135	ng/ul#	91
25) Dibenzo(a,h)anthracene	25.68	278	642	0.044	ng/ul#	1
26) Benzo(g,h,i)perylene	26.36	276	2228	0.144	ng/ul#	78

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

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