

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN101018\
 Data File : BN003059.D
 Acq On : 10 Oct 2018 10:10
 Operator : MJ/SJ
 Sample : J5115-13DL 5X
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 JLC46DL

Manual Integrations
 APPROVED

Sohil
 10/10/2018 4:58:32 PM

Quant Time: Oct 10 12:04:53 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 : Wed Oct 10 08:36:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1797	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.40	136	7695	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	4487	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	10425	0.40	ng/ul	-0.01
16) Chrysene-d12	21.22	240	9039	0.40	ng/ul	-0.01
20) Perylene-d12	23.44	264	7379	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	703	0.06	ng/ul	0.01
14) Fluoranthene-d10	19.05	212	1869	0.06	ng/ul	-0.01
Target Compounds				Qvalue		
5) 2-Methylnaphthalene	12.10	142	281	0.021	ng/ul	97
12) Phenanthrene	17.06	178	6624	0.228	ng/ul	98
13) Anthracene	17.15	178	1616	0.059	ng/ul#	89
15) Fluoranthene	19.09	202	25403	0.745	ng/ul	99
17) Pyrene	19.45	202	28303	0.806	ng/ul	97
18) Benzo(a)anthracene	21.21	228	14678	0.505	ng/ul	97
19) Chrysene	21.26	228	31630	1.071	ng/ul	99
21) Benzo(b)fluoranthene	22.77	252	30257m	1.139	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	8204m	0.298	ng/ul	
23) Benzo(a)pyrene	23.34	252	10608	0.445	ng/ul#	73
24) Indeno(1,2,3-cd)pyrene	25.65	276	11658	0.426	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.64	278	2860	0.131	ng/ul#	42
26) Benzo(g,h,i)perylene	26.32	276	12232	0.530	ng/ul	97

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

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