

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121918\
 Data File : BN004109.D
 Acq On : 19 Dec 2018 12:36
 Operator : JU/SJ
 Sample : J6297-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E37

Manual Integrations
 APPROVED

Sohil
 12/20/2018 2:02:52 PM

Quant Time: Dec 20 05:02:53 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 20 04:42:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	1242	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	6351	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	4070	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	10960m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	11893m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	10525m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	2347	0.24	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	7742m	0.24	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.19	152	1083	0.059	ng/ul#	85
12) Phenanthrene	17.25	178	1207m	0.034	ng/ul	
13) Anthracene	17.35	178	1201m	0.041	ng/ul	
15) Fluoranthene	19.27	202	3654m	0.086	ng/ul	
17) Pyrene	19.63	202	4600m	0.108	ng/ul	
18) Benzo(a)anthracene	21.39	228	3298	0.088	ng/ul	93
19) Chrysene	21.44	228	3879m	0.090	ng/ul	
21) Benzo(b)fluoranthene	23.02	252	6763m	0.155	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	2449m	0.058	ng/ul	
23) Benzo(a)pyrene	23.62	252	2656	0.070	ng/ul#	71
24) Indeno(1,2,3-cd)pyrene	26.08	276	2899	0.067	ng/ul#	97
26) Benzo(g,h,i)perylene	26.81	276	2348	0.062	ng/ul#	77

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

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Quant Time: Dec 20 05:02:53 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN11918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
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Manual Integrations
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