

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082318\
 Data File : BN002475.D
 Acq On : 23 Aug 2018 20:39
 Operator : JU/SJ
 Sample : J4421-11
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AA9

Manual Integrations
 APPROVED

Sohil
 8/24/2018 5:11:47 PM

Quant Time: Aug 24 01:58:54 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN082318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 24 00:52:48 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	1539	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	6382	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.31	164	3749	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	9007m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	9559m	0.40	ng/ul	0.00
20) Perylene-d12	23.48	264	8758m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	2293	0.23	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	6294m	0.24	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.09	178	551m	0.021	ng/ul	
15) Fluoranthene	19.12	202	1030m	0.033	ng/ul	
17) Pyrene	19.48	202	1293	0.040	ng/ul#	83
18) Benzo(a)anthracene	21.25	228	814	0.028	ng/ul#	87
19) Chrysene	21.30	228	855	0.026	ng/ul#	89
21) Benzo(b)fluoranthene	22.82	252	1208	0.038	ng/ul#	51
24) Indeno(1,2,3-cd)pyrene	25.71	276	797	0.026	ng/ul#	81
26) Benzo(g,h,i)perylene	26.39	276	758	0.029	ng/ul#	75

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

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