

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121518\
 Data File : BN003970.D
 Acq On : 15 Dec 2018 22:10
 Operator : JU/SJ
 Sample : J6171-06
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9Y10

Manual Integrations
 APPROVED

Sohil
 12/17/2018 4:24:00 PM

Quant Time: Dec 15 23:25:15 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 : Sat Dec 15 23:04:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	1887	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	9506	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5714	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	14447	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	18022	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	15906m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	2118	0.14	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	6414	0.15	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.68	128	1291	0.048	ng/ul#	83
5) 2-Methylnaphthalene	12.29	142	2262	0.121	ng/ul	100
7) Acenaphthylene	14.19	152	1927	0.075	ng/ul#	88
12) Phenanthrene	17.26	178	9494	0.200	ng/ul	97
13) Anthracene	17.35	178	2820	0.072	ng/ul#	90
15) Fluoranthene	19.27	202	14637	0.261	ng/ul	94
17) Pyrene	19.64	202	18882m	0.292	ng/ul	
18) Benzo(a)anthracene	21.38	228	10096	0.177	ng/ul#	84
19) Chrysene	21.44	228	16698	0.256	ng/ul#	90
21) Benzo(b)fluoranthene	23.01	252	20056m	0.305	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	5775m	0.090	ng/ul	
23) Benzo(a)pyrene	23.61	252	9599	0.168	ng/ul#	54
24) Indeno(1,2,3-cd)pyrene	26.07	276	9186	0.141	ng/ul#	93
25) Dibenzo(a,h)anthracene	26.08	278	2612	0.050	ng/ul#	1
26) Benzo(g,h,i)perylene	26.80	276	11518	0.203	ng/ul#	80

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

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