

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121518\
 Data File : BN003965.D
 Acq On : 15 Dec 2018 19:14
 Operator : JU/SJ
 Sample : J6227-20
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E69

Manual Integrations
 APPROVED

Sohil
 12/17/2018 4:23:55 PM

Quant Time: Dec 15 23:12:29 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	2173	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	10579	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	6313	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	15409m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	18733m	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	16015m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	3077	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	9125m	0.20	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.68	128	1227	0.041	ng/ul#	83
5) 2-Methylnaphthalene	12.29	142	2511	0.121	ng/ul	99
12) Phenanthrene	17.26	178	13430	0.266	ng/ul	99
13) Anthracene	17.35	178	2142m	0.052	ng/ul	
15) Fluoranthene	19.27	202	4049m	0.068	ng/ul	
17) Pyrene	19.63	202	15733m	0.234	ng/ul	
18) Benzo(a)anthracene	21.38	228	7616	0.129	ng/ul#	88
19) Chrysene	21.43	228	13438	0.198	ng/ul#	95
21) Benzo(b)fluoranthene	23.01	252	4439m	0.067	ng/ul	
22) Benzo(k)fluoranthene	23.05	252	1603m	0.025	ng/ul	
23) Benzo(a)pyrene	23.61	252	4255	0.074	ng/ul#	68
24) Indeno(1,2,3-cd)pyrene	26.06	276	2618	0.040	ng/ul#	99
25) Dibenzo(a,h)anthracene	26.07	278	1720	0.032	ng/ul#	20
26) Benzo(g,h,i)perylene	26.79	276	4578	0.080	ng/ul#	80

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

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