

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061819\
 Data File : BN006384.D
 Acq On : 18 Jun 2019 19:34
 Operator : JU/SJ
 Sample : K3335-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4224

Manual Integrations
 APPROVED

mohammad
 6/21/2019 8:39:20 AM

Quant Time: Jun 19 01:58:57 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 17 11:13:07 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	4024	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	15217	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	7783	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	14862m	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	11316m	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	10620m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	5521	0.24	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	11348m	0.28	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.01	152	1771	0.041	ng/ul#	89
12) Phenanthrene	17.09	178	14413	0.302	ng/ul	98
13) Anthracene	17.18	178	3487m	0.077	ng/ul	
15) Fluoranthene	19.11	202	37819m	0.718	ng/ul	
17) Pyrene	19.48	202	33643m	0.564	ng/ul	
18) Benzo(a)anthracene	21.25	228	14937	0.298	ng/ul	98
19) Chrysene	21.30	228	14865	0.316	ng/ul	98
21) Benzo(b)fluoranthene	22.83	252	20843m	0.471	ng/ul	
22) Benzo(k)fluoranthene	22.87	252	7531m	0.175	ng/ul	
23) Benzo(a)pyrene	23.40	252	13828m	0.335	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.73	276	9710m	0.218	ng/ul	
25) Dibenzo(a,h)anthracene	25.73	278	2011m	0.056	ng/ul	
26) Benzo(g,h,i)perylene	26.41	276	8224m	0.220	ng/ul	

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

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