

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN061419\
 Data File : BN006330.D
 Acq On : 14 Jun 2019 11:38
 Operator : JU/SJ
 Sample : K3332-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4226

Manual Integrations
 APPROVED

mohammad
 6/17/2019 8:28:43 AM

Quant Time: Jun 14 16:52:12 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 : Thu Jun 13 11:20:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	4414	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	16830	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	8269	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	14854m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	10084m	0.40	ng/ul	0.00
20) Perylene-d12	23.50	264	9705m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	5158	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	9712m	0.24	ng/ul	0.00
Target Compounds				Qvalue		
7) Acenaphthylene	14.01	152	1861	0.040	ng/ul#	89
8) Acenaphthene	14.36	153	643	0.020	ng/ul	95
9) Fluorene	15.35	166	1058	0.028	ng/ul#	94
12) Phenanthrene	17.09	178	18539	0.389	ng/ul	99
13) Anthracene	17.18	178	4882m	0.107	ng/ul	
15) Fluoranthene	19.12	202	36459m	0.692	ng/ul	
17) Pyrene	19.49	202	31223	0.588	ng/ul	99
18) Benzo(a)anthracene	21.25	228	16214	0.364	ng/ul	99
19) Chrysene	21.30	228	15246	0.364	ng/ul	98
21) Benzo(b)fluoranthene	22.83	252	21419m	0.529	ng/ul	
22) Benzo(k)fluoranthene	22.87	252	8731m	0.222	ng/ul	
23) Benzo(a)pyrene	23.40	252	13389m	0.355	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.73	276	9960m	0.244	ng/ul	
25) Dibenzo(a,h)anthracene	25.73	278	2854m	0.088	ng/ul	
26) Benzo(g,h,i)perylene	26.42	276	7923m	0.232	ng/ul	

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

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