

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN061419\
 Data File : BN006334.D
 Acq On : 14 Jun 2019 13:56
 Operator : JU/SJ
 Sample : K3332-05
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4230

Manual Integrations
 APPROVED

mohammad
 6/17/2019 8:28:52 AM

Quant Time: Jun 14 17:32:06 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	4002	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	15040	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	7640	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	15136m	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	11584m	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	10769m	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	3548	0.15	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	9921m	0.24	ng/ul	0.00
Target Compounds				Qvalue		
7) Acenaphthylene	14.01	152	4278	0.101	ng/ul	97
12) Phenanthrene	17.09	178	6084	0.125	ng/ul	96
13) Anthracene	17.18	178	4249m	0.092	ng/ul	
15) Fluoranthene	19.12	202	28018m	0.522	ng/ul	
17) Pyrene	19.49	202	28617	0.469	ng/ul	99
18) Benzo(a)anthracene	21.26	228	23481	0.458	ng/ul	99
19) Chrysene	21.31	228	21823	0.454	ng/ul	99
21) Benzo(b)fluoranthene	22.85	252	32019m	0.713	ng/ul	
22) Benzo(k)fluoranthene	22.88	252	13229m	0.302	ng/ul	
23) Benzo(a)pyrene	23.42	252	17052m	0.408	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.76	276	11670m	0.258	ng/ul	
25) Dibenzo(a,h)anthracene	25.76	278	3092m	0.085	ng/ul	
26) Benzo(g,h,i)perylene	26.44	276	8664m	0.229	ng/ul	

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

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