

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060819\
 Data File : BN006203.D
 Acq On : 08 Jun 2019 23:06
 Operator : JU/SJ
 Sample : K3016-04
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A51C3

Manual Integrations
 APPROVED

mohammad
 6/11/2019 4:31:59 PM

Quant Time: Jun 09 05:54:51 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 : Sun Jun 09 05:10:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	4654	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	17643	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	9633	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	20630m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	16440m	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	16765m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	7334	0.27	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	18316m	0.33	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.47	128	9820	0.202	ng/ul	96
5) 2-Methylnaphthalene	12.10	142	4763	0.140	ng/ul	97
7) Acenaphthylene	14.01	152	101004	1.886	ng/ul	100
8) Acenaphthene	14.36	153	4455	0.120	ng/ul	97
9) Fluorene	15.35	166	14785	0.340	ng/ul#	93
12) Phenanthrene	17.09	178	236333m	3.569	ng/ul	
13) Anthracene	17.18	178	89711m	1.419	ng/ul	
15) Fluoranthene	19.12	202	552376m	7.554	ng/ul	
17) Pyrene	19.48	202	590713	6.821	ng/ul	97
18) Benzo(a)anthracene	21.24	228	295566m	4.065	ng/ul	
19) Chrysene	21.29	228	341330	5.000	ng/ul	98
21) Benzo(b)fluoranthene	22.82	252	452773m	6.477	ng/ul	
22) Benzo(k)fluoranthene	22.86	252	171490m	2.519	ng/ul	
23) Benzo(a)pyrene	23.39	252	312605	4.801	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	25.71	276	225073	3.194	ng/ul#	89
25) Dibenzo(a,h)anthracene	25.72	278	57686	1.024	ng/ul	100
26) Benzo(a,h,i)perylene	26.40	276	173013	2.931	ng/ul	95

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

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