

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120619\
 Data File : BN008866.D
 Acq On : 06 Dec 2019 19:02
 Operator : JU
 Sample : K6007-12
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BL8

Manual Integrations
 APPROVED

mohammad
 12/9/2019 3:07:45 PM

Quant Time: Dec 06 23:11:08 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 06 12:03:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	3017	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	13766	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	9088	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	20421	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	18714	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	21228	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	3649	0.16	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	11028	0.18	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.37	128	2913	0.070	ng/ul#	95
5) 2-Methylnaphthalene	12.00	142	2172	0.075	ng/ul	98
7) Acenaphthylene	13.91	152	5025	0.100	ng/ul	97
9) Fluorene	15.25	166	1143	0.026	ng/ul#	92
12) Phenanthrene	16.98	178	16860	0.245	ng/ul	100
13) Anthracene	17.07	178	4908	0.075	ng/ul#	95
15) Fluoranthene	19.00	202	56331	0.666	ng/ul	99
17) Pyrene	19.37	202	48523	0.646	ng/ul	99
18) Benzo(a)anthracene	21.12	228	34220	0.455	ng/ul	97
19) Chrysene	21.17	228	33118	0.448	ng/ul	97
21) Benzo(b)fluoranthene	22.66	252	47600m	0.544	ng/ul	
22) Benzo(k)fluoranthene	22.69	252	14735m	0.170	ng/ul	
23) Benzo(a)pyrene	23.20	252	26984	0.328	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.40	276	17773	0.185	ng/ul	93
25) Dibenzo(a,h)anthracene	25.40	278	5045	0.065	ng/ul#	78
26) Benzo(a,h,i)perylene	26.04	276	15321	0.191	ng/ul	99

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

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