

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120619\
 Data File : BN008857.D
 Acq On : 06 Dec 2019 13:40
 Operator : JU
 Sample : K6007-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BK9

Manual Integrations
 APPROVED

mohammad
 12/9/2019 3:06:56 PM

Quant Time: Dec 06 14:54:53 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	2514	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	11379	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	7468	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	16579m	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	14818	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	16725	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	4953	0.27	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	13926	0.28	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.37	128	3742	0.109	ng/ul#	94
5) 2-Methylnaphthalene	12.00	142	2780	0.116	ng/ul	100
7) Acenaphthylene	13.91	152	1478	0.036	ng/ul#	85
9) Fluorene	15.25	166	1517	0.042	ng/ul#	91
12) Phenanthrene	16.98	178	19198	0.343	ng/ul	100
13) Anthracene	17.07	178	2185	0.041	ng/ul#	91
15) Fluoranthene	19.00	202	24867	0.362	ng/ul	100
17) Pyrene	19.37	202	24296	0.409	ng/ul	98
18) Benzo(a)anthracene	21.13	228	11622	0.195	ng/ul#	91
19) Chrysene	21.18	228	17210	0.294	ng/ul#	93
21) Benzo(b)fluoranthene	22.66	252	17890m	0.259	ng/ul	
22) Benzo(k)fluoranthene	22.70	252	7188m	0.105	ng/ul	
23) Benzo(a)pyrene	23.20	252	11871	0.183	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.41	276	9290	0.123	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.42	278	2607	0.043	ng/ul#	47
26) Benzo(a,h,i)perylene	26.05	276	9965	0.157	ng/ul	96

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

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