

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120619\
 Data File : BN008867.D
 Acq On : 06 Dec 2019 19:37
 Operator : JU
 Sample : K6007-13
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BL9

Manual Integrations
 APPROVED

mohammad
 12/9/2019 3:07:49 PM

Quant Time: Dec 06 23:13:35 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	7.56	152	2915	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	13230	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	8889	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	19705	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	17832	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	19487	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	2967	0.14	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	9352	0.16	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.37	128	3981	0.100	ng/ul	96
5) 2-Methylnaphthalene	12.00	142	3448	0.124	ng/ul	99
7) Acenaphthylene	13.91	152	24588	0.500	ng/ul	99
8) Acenaphthene	14.25	153	809	0.022	ng/ul#	94
9) Fluorene	15.25	166	2463	0.058	ng/ul#	90
12) Phenanthrene	16.98	178	73542	1.107	ng/ul	99
13) Anthracene	17.07	178	21019	0.335	ng/ul	99
15) Fluoranthene	19.00	202	230117	2.819	ng/ul	98
17) Pyrene	19.37	202	174989	2.446	ng/ul	98
18) Benzo(a)anthracene	21.12	228	137517	1.918	ng/ul	95
19) Chrysene	21.18	228	132791	1.885	ng/ul	98
21) Benzo(b)fluoranthene	22.66	252	181939m	2.264	ng/ul	
22) Benzo(k)fluoranthene	22.69	252	60229m	0.756	ng/ul	
23) Benzo(a)pyrene	23.20	252	104945	1.391	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	25.40	276	75155	0.852	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.41	278	22167	0.311	ng/ul	95
26) Benzo(a,h,i)perylene	26.05	276	63824	0.865	ng/ul	97

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

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