

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102419\
 Data File : BN008415.D
 Acq On : 25 Oct 2019 07:31
 Operator : JU
 Sample : K5236-10
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPL9

Manual Integrations
 APPROVED

mohammad
 10/25/2019 9:28:24 PM

Quant Time: Oct 25 12:06:18 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 25 00:45:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	1658	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	8381	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.21	164	5400	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	12279	0.40	ng/ul	0.00
16) Chrysene-d12	21.17	240	11116	0.40	ng/ul	0.00
20) Perylene-d12	23.34	264	9108	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	2242	0.16	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	7588	0.18	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.38	128	5998	0.251	ng/ul	99
5) 2-Methylnaphthalene	12.01	142	2285	0.140	ng/ul	99
7) Acenaphthylene	13.93	152	36806	1.431	ng/ul	99
8) Acenaphthene	14.27	153	1981	0.099	ng/ul	95
9) Fluorene	15.26	166	15149	0.667	ng/ul	99
12) Phenanthrene	17.00	178	252683	7.029	ng/ul	99
13) Anthracene	17.09	178	46621	1.474	ng/ul	99
15) Fluoranthene	19.03	202	450552	9.489	ng/ul	99
17) Pyrene	19.39	202	320799	7.248	ng/ul	98
18) Benzo(a)anthracene	21.15	228	185200	4.583	ng/ul	97
19) Chrysene	21.20	228	175131	3.531	ng/ul	97
21) Benzo(b)fluoranthene	22.70	252	196637m	6.632	ng/ul	
22) Benzo(k)fluoranthene	22.73	252	76144m	2.267	ng/ul	
23) Benzo(a)pyrene	23.25	252	120500	3.862	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.49	276	92348	2.513	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.49	278	21936	0.755	ng/ul#	89
26) Benzo(a,h,i)perylene	26.14	276	102089	3.030	ng/ul	98

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

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