

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102219\
 Data File : BN008331.D
 Acq On : 23 Oct 2019 00:57
 Operator : JU
 Sample : K5223-08
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPE4

Manual Integrations
 APPROVED

mohammad
 10/24/2019 8:05:49 AM

Quant Time: Oct 23 02:46:44 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 : Mon Oct 21 23:58:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	1940	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.35	136	9470	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.22	164	5921	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.97	188	13047m	0.40	ng/ul	0.00
16) Chrysene-d12	21.17	240	11394	0.40	ng/ul	0.00
20) Perylene-d12	23.34	264	10212	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	3843	0.24	ng/ul	-0.01
14) Fluoranthene-d10	19.00	212	10438	0.23	ng/ul	0.00
Target Compounds				Ovalue		
8) Acenaphthene	14.28	153	467	0.021	ng/ul	95
9) Fluorene	15.27	166	534	0.021	ng/ul#	93
12) Phenanthrene	17.01	178	6588	0.172	ng/ul	98
13) Anthracene	17.10	178	1314	0.039	ng/ul#	87
15) Fluoranthene	19.03	202	6656	0.132	ng/ul	95
17) Pyrene	19.39	202	5782m	0.127	ng/ul	
18) Benzo(a)anthracene	21.16	228	3238	0.078	ng/ul#	90
19) Chrysene	21.21	228	2940	0.058	ng/ul#	92
21) Benzo(b)fluoranthene	22.70	252	2563m	0.077	ng/ul	
22) Benzo(k)fluoranthene	22.74	252	1066m	0.028	ng/ul	
23) Benzo(a)pyrene	23.25	252	1788	0.051	ng/ul#	23
26) Benzo(g,h,i)perylene	26.16	276	799	0.021	ng/ul#	46

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

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