

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102119\
 Data File : BN008276.D
 Acq On : 21 Oct 2019 11:01
 Operator : JU
 Sample : K5199-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPJ8

Manual Integrations
 APPROVED

mohammad
 10/22/2019 3:26:56 PM

Quant Time: Oct 21 15:41:05 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 15:18:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	1784	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	8529	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	5338	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.97	188	12124m	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	11800	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	13916	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	3920	0.27	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	11613	0.28	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.94	152	575	0.023	ng/ul#	68
9) Fluorene	15.28	166	472	0.021	ng/ul#	94
12) Phenanthrene	17.01	178	10985	0.309	ng/ul	99
13) Anthracene	17.11	178	1879	0.060	ng/ul#	89
15) Fluoranthene	19.04	202	21429	0.457	ng/ul	96
17) Pyrene	19.40	202	18386	0.391	ng/ul	98
18) Benzo(a)anthracene	21.16	228	9108	0.212	ng/ul	95
19) Chrysene	21.22	228	9883	0.188	ng/ul	97
21) Benzo(b)fluoranthene	22.71	252	12891m	0.285	ng/ul	
22) Benzo(k)fluoranthene	22.75	252	4613m	0.090	ng/ul	
23) Benzo(a)pyrene	23.26	252	7494	0.157	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.51	276	4306	0.077	ng/ul#	90
25) Dibenzo(a,h)anthracene	25.51	278	941	0.021	ng/ul#	28
26) Benzo(g,h,i)perylene	26.16	276	3593	0.070	ng/ul#	84

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

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