

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102219\
 Data File : BN008323.D
 Acq On : 22 Oct 2019 20:10
 Operator : JU
 Sample : K5223-01
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPH3

Manual Integrations
 APPROVED

mohammad
 10/24/2019 8:05:27 AM

Quant Time: Oct 23 02:10:19 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	1657	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	8241	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.22	164	5205	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.97	188	11330m	0.40	ng/ul	0.00
16) Chrysene-d12	21.17	240	10795	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	9693	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	3116	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	9854	0.25	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.39	128	547	0.023	ng/ul#	68
7) Acenaphthylene	13.94	152	1384	0.056	ng/ul#	81
9) Fluorene	15.27	166	501	0.023	ng/ul#	91
12) Phenanthrene	17.01	178	10363	0.312	ng/ul	100
13) Anthracene	17.10	178	2280	0.078	ng/ul#	93
15) Fluoranthene	19.03	202	26393	0.602	ng/ul	96
17) Pyrene	19.40	202	23473m	0.546	ng/ul	
18) Benzo(a)anthracene	21.16	228	15082	0.384	ng/ul	96
19) Chrysene	21.21	228	16935	0.352	ng/ul	97
21) Benzo(b)fluoranthene	22.70	252	18453m	0.585	ng/ul	
22) Benzo(k)fluoranthene	22.74	252	6272m	0.175	ng/ul	
23) Benzo(a)pyrene	23.25	252	12349	0.372	ng/ul#	95
24) Indeno(1,2,3-cd)pyrene	25.50	276	7767	0.199	ng/ul#	90
25) Dibenzo(a,h)anthracene	25.50	278	2179	0.070	ng/ul#	64
26) Benzo(g,h,i)perylene	26.15	276	5648	0.157	ng/ul#	90

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

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