

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012721\
 Data File : BN013519.D
 Acq On : 27 Jan 2021 23:07
 Operator : CG/JU
 Sample : M1099-03
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F1C03

Manual Integrations
 APPROVED

mohammad
 1/28/2021 11:44:45 AM

Quant Time: Jan 28 06:55:22 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN012721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 28 06:42:07 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	4387	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	16594	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	9601	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	19692m	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	18526	0.40	ng/ul	0.00
20) Perylene-d12	23.32	264	18616	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	10455	0.42	ng/ul	0.00
14) Fluoranthene-d10	18.99	212	21238	0.41	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.38	128	1765	0.037	ng/ul#	89
7) Acenaphthylene	13.92	152	1943	0.049	ng/ul#	90
9) Fluorene	15.25	166	785	0.020	ng/ul#	90
12) Phenanthrene	16.99	178	4472	0.069	ng/ul	96
15) Fluoranthene	19.01	202	7621	0.108	ng/ul	97
17) Pyrene	19.38	202	11937	0.157	ng/ul	98
18) Benzo(a)anthracene	21.13	228	2935m	0.048	ng/ul	
19) Chrysene	21.18	228	3689	0.048	ng/ul	94
21) Benzo(b)fluoranthene	22.67	252	5973m	0.076	ng/ul	
22) Benzo(k)fluoranthene	22.71	252	2129m	0.026	ng/ul	
23) Benzo(a)pyrene	23.23	252	3687	0.056	ng/ul#	56
24) Indeno(1,2,3-cd)pyrene	25.47	276	3110	0.035	ng/ul#	100
26) Benzo(g,h,i)perylene	26.13	276	3649	0.047	ng/ul#	89

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

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