

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012721\
 Data File : BN013516.D
 Acq On : 27 Jan 2021 21:20
 Operator : CG/JU
 Sample : M1099-01
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F1C01

Manual Integrations
 APPROVED

mohammad
 1/28/2021 11:44:31 AM

Quant Time: Jan 28 06:51:02 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	5008	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	19364	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	11537	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	24107m	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	22287	0.40	ng/ul	0.00
20) Perylene-d12	23.32	264	21487	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	10600	0.37	ng/ul	0.00
14) Fluoranthene-d10	18.99	212	23094	0.36	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.38	128	3784	0.068	ng/ul#	94
5) 2-Methylnaphthalene	12.00	142	822	0.023	ng/ul	95
7) Acenaphthylene	13.92	152	4273	0.090	ng/ul	94
8) Acenaphthene	14.26	153	2917	0.069	ng/ul	97
9) Fluorene	15.25	166	3223	0.069	ng/ul#	97
12) Phenanthrene	16.99	178	22413	0.284	ng/ul	98
13) Anthracene	17.08	178	7789	0.120	ng/ul	97
15) Fluoranthene	19.01	202	18062	0.209	ng/ul	98
17) Pyrene	19.38	202	17795	0.195	ng/ul	99
18) Benzo(a)anthracene	21.13	228	15130	0.206	ng/ul	95
19) Chrysene	21.19	228	24492	0.267	ng/ul	99
21) Benzo(b)fluoranthene	22.67	252	23012m	0.255	ng/ul	
22) Benzo(k)fluoranthene	22.71	252	10216m	0.106	ng/ul	
23) Benzo(a)pyrene	23.23	252	15635	0.204	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.47	276	8729	0.086	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.48	278	2338	0.028	ng/ul#	63
26) Benzo(a,h,i)perylene	26.13	276	8705	0.097	ng/ul	96

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

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