

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012721\
 Data File : BN013505.D
 Acq On : 27 Jan 2021 13:34
 Operator : CG/JU
 Sample : M1099-16
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F1C12

Manual Integrations
 APPROVED

mohammad
 1/28/2021 11:44:10 AM

Quant Time: Jan 27 14:06:24 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 : Wed Jan 27 11:55:49 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	5322	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	20585	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	12285	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	25109m	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	23262	0.40	ng/ul	0.00
20) Perylene-d12	23.32	264	22440	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	10544	0.34	ng/ul	0.00
14) Fluoranthene-d10	18.99	212	22923	0.34	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.38	128	20985	0.352	ng/ul	98
5) 2-Methylnaphthalene	12.00	142	6955	0.184	ng/ul	99
7) Acenaphthylene	13.92	152	65746	1.299	ng/ul	100
8) Acenaphthene	14.26	153	23194	0.515	ng/ul	99
9) Fluorene	15.25	166	27442	0.551	ng/ul	98
11) Pentachlorophenol	16.60	266	76	0.021	ng/ul	93
12) Phenanthrene	16.99	178	111090	1.351	ng/ul	99
13) Anthracene	17.08	178	21185	0.314	ng/ul	98
15) Fluoranthene	19.01	202	108873	1.212	ng/ul	99
17) Pyrene	19.38	202	213400	2.241	ng/ul	99
18) Benzo(a)anthracene	21.14	228	10001	0.130	ng/ul#	1
19) Chrysene	21.19	228	13331	0.139	ng/ul#	89
21) Benzo(b)fluoranthene	22.68	252	13015m	0.138	ng/ul	
22) Benzo(k)fluoranthene	22.72	252	4095m	0.041	ng/ul	
23) Benzo(a)pyrene	23.23	252	13235	0.166	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	25.48	276	11228	0.106	ng/ul#	100
26) Benzo(a,h,i)perylene	26.13	276	30358	0.323	ng/ul	99

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

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