

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120720\
 Data File : BN012954.D
 Acq On : 08 Dec 2020 00:08
 Operator : CG/JU
 Sample : L4862-14
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0B04

Manual Integrations
 APPROVED

mohammad
 12/9/2020 12:21:51 PM

Quant Time: Dec 08 02:06:51 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN120320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 08 00:15:39 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1028	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	4249	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.05	164	2545	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	5324m	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	4620	0.40	ng/ul	0.00
20) Perylene-d12	23.13	264	4518	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.75	152	2792	0.44	ng/ul	0.00
14) Fluoranthene-d10	18.85	212	6116	0.42	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.20	128	284	0.023	ng/ul#	69
5) 2-Methylnaphthalene	11.83	142	187	0.023	ng/ul	93
7) Acenaphthylene	13.76	152	280	0.023	ng/ul#	57
12) Phenanthrene	16.85	178	3141	0.176	ng/ul	99
13) Anthracene	16.94	178	548	0.036	ng/ul#	85
15) Fluoranthene	18.88	202	11639	0.549	ng/ul	99
17) Pyrene	19.25	202	12016	0.551	ng/ul	97
18) Benzo(a)anthracene	21.01	228	6682	0.350	ng/ul	99
19) Chrysene	21.07	228	8374	0.397	ng/ul	99
21) Benzo(b)fluoranthene	22.51	252	12558m	0.601	ng/ul	
22) Benzo(k)fluoranthene	22.55	252	4843m	0.210	ng/ul	
23) Benzo(a)pyrene	23.04	252	7765	0.400	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.19	276	5311	0.208	ng/ul#	90
25) Dibenzo(a,h)anthracene	25.19	278	1335	0.066	ng/ul#	66
26) Benzo(g,h,i)perylene	25.82	276	5363	0.233	ng/ul	96

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

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