

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120720\
 Data File : BN012967.D
 Acq On : 08 Dec 2020 09:09
 Operator : CG/JU
 Sample : L4859-04
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0B94

Manual Integrations
 APPROVED

mohammad
 12/9/2020 12:24:26 PM

Quant Time: Dec 08 12:04:29 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 : Mon Dec 07 12:24:06 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1168	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	4619	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.05	164	2499	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	5236	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	4192	0.40	ng/ul	0.00
20) Perylene-d12	23.13	264	3686	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.75	152	736	0.11	ng/ul	0.00
14) Fluoranthene-d10	18.85	212	1643	0.12	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.20	128	1075	0.080	ng/ul	95
5) 2-Methylnaphthalene	11.83	142	551	0.063	ng/ul	100
7) Acenaphthylene	13.76	152	3478	0.294	ng/ul	98
8) Acenaphthene	14.11	153	1948	0.201	ng/ul	99
9) Fluorene	15.11	166	2439	0.219	ng/ul	99
12) Phenanthrene	16.85	178	66521	3.789	ng/ul	97
13) Anthracene	16.94	178	14827	0.992	ng/ul	95
15) Fluoranthene	18.88	202	211724	10.149	ng/ul	97
17) Pyrene	19.25	202	180271	9.104	ng/ul	99
18) Benzo(a)anthracene	21.02	228	104321	6.015	ng/ul	99
19) Chrysene	21.07	228	108547	5.664	ng/ul	98
21) Benzo(b)fluoranthene	22.52	252	135059m	7.920	ng/ul	
22) Benzo(k)fluoranthene	22.56	252	46647m	2.475	ng/ul	
23) Benzo(a)pyrene	23.04	252	85543	5.406	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	25.19	276	57252	2.755	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.20	278	14176	0.862	ng/ul	94
26) Benzo(a,h,i)perylene	25.82	276	48486	2.579	ng/ul	97

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

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