

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120620\
 Data File : BN012928.D
 Acq On : 06 Dec 2020 21:30
 Operator : CG/JU
 Sample : L4864-20
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0B63

Manual Integrations
 APPROVED

mohammad
 12/7/2020 2:12:05 PM

Quant Time: Dec 07 04:05:29 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 03:06:53 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1420	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	5830	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.06	164	3298	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	6872	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	5847	0.40	ng/ul	0.00
20) Perylene-d12	23.14	264	5730	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	946	0.11	ng/ul	-0.02
14) Fluoranthene-d10	18.86	212	2134	0.11	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.20	128	454	0.027	ng/ul#	79
5) 2-Methylnaphthalene	11.84	142	271	0.024	ng/ul	100
7) Acenaphthylene	13.77	152	435	0.028	ng/ul#	73
8) Acenaphthene	14.12	153	564	0.044	ng/ul	97
9) Fluorene	15.12	166	729	0.050	ng/ul#	98
12) Phenanthrene	16.85	178	24476	1.062	ng/ul	97
13) Anthracene	16.95	178	3849	0.196	ng/ul	98
15) Fluoranthene	18.88	202	73527	2.685	ng/ul	99
17) Pyrene	19.25	202	62283	2.255	ng/ul	98
18) Benzo(a)anthracene	21.02	228	30026	1.241	ng/ul	98
19) Chrysene	21.07	228	38505	1.441	ng/ul	98
21) Benzo(b)fluoranthene	22.52	252	54401m	2.052	ng/ul	
22) Benzo(k)fluoranthene	22.55	252	17509m	0.598	ng/ul	
23) Benzo(a)pyrene	23.04	252	30992m	1.260	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.19	276	21832	0.676	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.20	278	5220	0.204	ng/ul	94
26) Benzo(a,h,i)perylene	25.82	276	19629	0.672	ng/ul	97

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

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