

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112020\
 Data File : BN012669.D
 Acq On : 20 Nov 2020 21:15
 Operator : CG/JU
 Sample : L4710-06
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AX2

Manual Integrations
 APPROVED

mohammad
 11/24/2020 8:56:18 AM

Quant Time: Nov 21 01:08:36 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN112020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 21 00:16:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.44	152	1021	0.40	ng/ul	0.00
2) Naphthalene-d8	10.21	136	4303	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.10	164	2458	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.85	188	5010	0.40	ng/ul	0.00
16) Chrysene-d12	21.07	240	4409	0.40	ng/ul	0.00
20) Perylene-d12	23.18	264	4351	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.81	152	1233	0.20	ng/ul	-0.01
14) Fluoranthene-d10	18.89	212	2559	0.20	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.26	128	2685	0.207	ng/ul	100
5) 2-Methylnaphthalene	11.89	142	1552	0.188	ng/ul	100
7) Acenaphthylene	13.82	152	3544	0.296	ng/ul	97
8) Acenaphthene	14.16	153	17672	1.749	ng/ul	98
9) Fluorene	15.16	166	16957	1.527	ng/ul	99
12) Phenanthrene	16.89	178	582824	33.433	ng/ul	95
13) Anthracene	16.98	178	65119	4.297	ng/ul	95
15) Fluoranthene	18.93	202	1174178	58.340	ng/ul	99
17) Pyrene	19.29	202	930632	47.188	ng/ul	95
18) Benzo(a)anthracene	21.05	228	417387	26.418	ng/ul	98
19) Chrysene	21.10	228	463852	23.503	ng/ul	98
21) Benzo(b)fluoranthene	22.56	252	647573m	32.353	ng/ul	
22) Benzo(k)fluoranthene	22.59	252	229642m	9.949	ng/ul	
23) Benzo(a)pyrene	23.09	252	420537	22.324	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.27	276	327436	13.292	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.27	278	74072m	3.742	ng/ul	
26) Benzo(a,h,i)perylene	25.90	276	315177	14.216	ng/ul	96

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

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