

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112020\
 Data File : BN012678.D
 Acq On : 21 Nov 2020 10:01
 Operator : CG/JU
 Sample : L4710-12DL 2X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0B32DL

Manual Integrations
 APPROVED

mohammad
 11/24/2020 8:56:40 AM

Quant Time: Nov 23 01:14:41 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 : Mon Nov 23 00:58:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.44	152	1103	0.40	ng/ul	0.00
2) Naphthalene-d8	10.20	136	4579	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.10	164	2519	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.85	188	5209	0.40	ng/ul	0.00
16) Chrysene-d12	21.06	240	4533	0.40	ng/ul	0.00
20) Perylene-d12	23.18	264	4657	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.81	152	566	0.09	ng/ul	-0.01
14) Fluoranthene-d10	18.89	212	1389	0.10	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.25	128	772	0.056	ng/ul#	85
7) Acenaphthylene	13.81	152	290	0.024	ng/ul#	61
12) Phenanthrene	16.89	178	4440	0.245	ng/ul	99
13) Anthracene	16.98	178	664	0.042	ng/ul#	83
15) Fluoranthene	18.92	202	17598	0.841	ng/ul	99
17) Pyrene	19.28	202	15961	0.787	ng/ul#	95
18) Benzo(a)anthracene	21.05	228	7991	0.492	ng/ul	99
19) Chrysene	21.10	228	11640	0.574	ng/ul	99
21) Benzo(b)fluoranthene	22.56	252	17098m	0.798	ng/ul	
22) Benzo(k)fluoranthene	22.59	252	5961m	0.241	ng/ul	
23) Benzo(a)pyrene	23.09	252	10299m	0.511	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.26	276	8368	0.317	ng/ul	97
25) Dibenzo(a,h)anthracene	25.27	278	1886	0.089	ng/ul#	70
26) Benzo(g,h,i)perylene	25.89	276	8065	0.340	ng/ul	97

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

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