

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120720\
 Data File : BN013046.D
 Acq On : 10 Dec 2020 14:29
 Operator : CG/JU
 Sample : L4941-02DL 5X
 Misc :
 ALS Vial : 116 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AX3DL

Manual Integrations
 APPROVED

mohammad
 12/11/2020 1:19:22 PM

Quant Time: Dec 10 15:17:49 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 : Thu Dec 10 10:29:27 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	891	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	3612	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.04	164	1987	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.80	188	4227	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	3648	0.40	ng/ul	0.00
20) Perylene-d12	23.13	264	3603	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	150	0.03	ng/ul	0.01
14) Fluoranthene-d10	18.84	212	374	0.03	ng/ul	0.00
Target Compounds				Ovalue		
5) 2-Methylnaphthalene	11.84	142	138	0.020	ng/ul	100
7) Acenaphthylene	13.76	152	521	0.055	ng/ul#	79
8) Acenaphthene	14.11	153	2089	0.271	ng/ul	99
9) Fluorene	15.11	166	2706	0.306	ng/ul	98
12) Phenanthrene	16.84	178	74997	5.292	ng/ul	96
13) Anthracene	16.93	178	12393	1.027	ng/ul	95
15) Fluoranthene	18.87	202	180856	10.739	ng/ul	96
17) Pyrene	19.24	202	144454	8.383	ng/ul	99
18) Benzo(a)anthracene	21.01	228	68697	4.552	ng/ul	98
19) Chrysene	21.06	228	80656	4.837	ng/ul	98
21) Benzo(b)fluoranthene	22.51	252	98123m	5.887	ng/ul	
22) Benzo(k)fluoranthene	22.55	252	36193m	1.964	ng/ul	
23) Benzo(a)pyrene	23.04	252	63951m	4.135	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.17	276	41352	2.035	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.18	278	10229	0.637	ng/ul	95
26) Benzo(a,h,i)perylene	25.80	276	34637	1.885	ng/ul	96

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

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