

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120720\
 Data File : BN013017.D
 Acq On : 09 Dec 2020 18:20
 Operator : CG/JU
 Sample : L4941-03
 Misc :
 ALS Vial : 87 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AX8

Manual Integrations
 APPROVED

mohammad
 12/11/2020 1:17:52 PM

Quant Time: Dec 10 01:06:08 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 : Wed Dec 09 07:33:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	1008	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.14	136	3965	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.05	164	2202	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	4572m	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	3694	0.40	ng/ul	0.00
20) Perylene-d12	23.14	264	3438	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.75	152	1162	0.20	ng/ul	-0.01
14) Fluoranthene-d10	18.85	212	2600	0.21	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.76	152	386	0.037	ng/ul#	62
9) Fluorene	15.11	166	218	0.022	ng/ul#	89
12) Phenanthrene	16.85	178	6592	0.430	ng/ul	99
13) Anthracene	16.94	178	906	0.069	ng/ul#	88
15) Fluoranthene	18.88	202	22051	1.211	ng/ul	99
17) Pyrene	19.25	202	18926	1.085	ng/ul	96
18) Benzo(a)anthracene	21.02	228	10410	0.681	ng/ul	98
19) Chrysene	21.07	228	12944	0.767	ng/ul	99
21) Benzo(b)fluoranthene	22.52	252	16723m	1.051	ng/ul	
22) Benzo(k)fluoranthene	22.55	252	6150m	0.350	ng/ul	
23) Benzo(a)pyrene	23.04	252	10101	0.684	ng/ul	92
24) Indeno(1,2,3-cd)pyrene	25.19	276	6763	0.349	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.19	278	1627	0.106	ng/ul#	64
26) Benzo(g,h,i)perylene	25.82	276	5627	0.321	ng/ul#	94

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

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