

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120720\
 Data File : BN013007.D
 Acq On : 09 Dec 2020 11:26
 Operator : CG/JU
 Sample : L4942-04DL 5X
 Misc :
 ALS Vial : 77 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0B96DL

Manual Integrations
 APPROVED

mohammad
 12/9/2020 5:39:24 PM

Quant Time: Dec 09 11:58:59 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	1272	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	5198	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.05	164	2822	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	5641	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	4200	0.40	ng/ul	0.00
20) Perylene-d12	23.13	264	3804	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	216	0.03	ng/ul	0.00
14) Fluoranthene-d10	18.85	212	437	0.03	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.76	152	404	0.030	ng/ul#	75
9) Fluorene	15.11	166	278	0.022	ng/ul#	88
12) Phenanthrene	16.85	178	8479	0.448	ng/ul	98
13) Anthracene	16.94	178	1546	0.096	ng/ul	95
15) Fluoranthene	18.88	202	18728	0.833	ng/ul	99
17) Pyrene	19.25	202	16869	0.850	ng/ul	98
18) Benzo(a)anthracene	21.02	228	8082	0.465	ng/ul	99
19) Chrysene	21.07	228	8526	0.444	ng/ul	98
21) Benzo(b)fluoranthene	22.52	252	10488m	0.596	ng/ul	
22) Benzo(k)fluoranthene	22.55	252	3472m	0.178	ng/ul	
23) Benzo(a)pyrene	23.04	252	6765	0.414	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.19	276	4067	0.190	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.19	278	1043	0.061	ng/ul#	49
26) Benzo(g,h,i)perylene	25.82	276	3469	0.179	ng/ul#	91

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

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