

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121220\
 Data File : BN013096.D
 Acq On : 12 Dec 2020 17:58
 Operator : CG/JU
 Sample : L5063-04DL 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AD6DL

Manual Integrations
 APPROVED

mohammad
 12/15/2020 12:17:08 PM

Quant Time: Dec 14 01:26:53 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 : Mon Dec 14 00:25:18 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.36	152	952	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.13	136	3949	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.03	164	3979	0.40	ng/ul	-0.01
10) Phenanthrene-d10	16.79	188	4686	0.40	ng/ul	-0.01
16) Chrysene-d12	21.02	240	4426	0.40	ng/ul	0.00
20) Perylene-d12	23.12	264	3913	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.71	152	784	0.13	ng/ul	-0.03
14) Fluoranthene-d10	18.84	212	1138	0.09	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.17	128	31222	2.732	ng/ul	95
5) 2-Methylnaphthalene	11.80	142	159026	21.186	ng/ul	99
7) Acenaphthylene	13.75	152	8292m	0.440	ng/ul	
8) Acenaphthene	14.09	153	14188	0.918	ng/ul	93
9) Fluorene	15.09	166	26730	1.508	ng/ul#	88
12) Phenanthrene	16.84	178	343447	21.860	ng/ul	97
13) Anthracene	16.92	178	29765	2.226	ng/ul#	91
15) Fluoranthene	18.87	202	46540	2.493	ng/ul	92
17) Pyrene	19.24	202	254435	12.171	ng/ul	99
18) Benzo(a)anthracene	21.01	228	26813	1.464	ng/ul#	35
19) Chrysene	21.06	228	35502	1.755	ng/ul#	71
21) Benzo(b)fluoranthene	22.50	252	16475m	0.910	ng/ul	
22) Benzo(k)fluoranthene	22.54	252	1362m	0.068	ng/ul	
23) Benzo(a)pyrene	23.03	252	14801	0.881	ng/ul#	35
24) Indeno(1,2,3-cd)pyrene	25.17	276	5052	0.229	ng/ul#	88
25) Dibenzo(a,h)anthracene	25.17	278	2790	0.160	ng/ul#	1
26) Benzo(a,h,i)perylene	25.80	276	9517	0.477	ng/ul#	58

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

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