

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN011020\
 Data File : BN009352.D
 Acq On : 11 Jan 2020 02:13
 Operator : JU
 Sample : L1022-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 H0BP6

Manual Integrations
 APPROVED

mohammad
 1/13/2020 2:22:40 PM

Quant Time: Jan 11 02:55:20 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN011020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 10 16:45:39 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.48	152	1302	0.40	ng/ul	0.00
2) Naphthalene-d8	10.22	136	4646	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.11	164	2730	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	5900m	0.40	ng/ul	0.00
16) Chrysene-d12	21.07	240	4840	0.40	ng/ul	0.00
20) Perylene-d12	23.18	264	5271	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.84	152	1854	0.25	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	5707	0.32	ng/ul	0.00
Target Compounds				Ovalue		
11) Pentachlorophenol	16.54	266	66	0.041	ng/ul#	91
12) Phenanthrene	16.90	178	539	0.026	ng/ul#	79
13) Anthracene	16.99	178	440	0.024	ng/ul#	72
15) Fluoranthene	18.93	202	664	0.028	ng/ul	76
17) Pyrene	19.29	202	697	0.030	ng/ul#	68
18) Benzo(a)anthracene	21.05	228	666	0.031	ng/ul#	84
19) Chrysene	21.10	228	799	0.035	ng/ul#	89
21) Benzo(b)fluoranthene	22.56	252	805	0.034	ng/ul#	24
22) Benzo(k)fluoranthene	22.60	252	737	0.028	ng/ul#	8
23) Benzo(a)pyrene	23.09	252	557	0.024	ng/ul#	11
24) Indeno(1,2,3-cd)pyrene	25.28	276	721	0.028	ng/ul#	75
25) Dibenzo(a,h)anthracene	25.28	278	501	0.024	ng/ul#	12
26) Benzo(g,h,i)perylene	25.90	276	662	0.029	ng/ul#	65

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

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