

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120720\
 Data File : BN012996.D
 Acq On : 09 Dec 2020 04:05
 Operator : CG/JU
 Sample : L4942-02
 Misc :
 ALS Vial : 66 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0B78

Quant Time: Dec 09 07:42:10 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	1030	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	4350	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.05	164	2488	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	5187	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	4081	0.40	ng/ul	0.00
20) Perylene-d12	23.13	264	3816	0.40	ng/ul	-0.01

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.75	152	1144	0.18	ng/ul	0.00
14) Fluoranthene-d10	18.85	212	2678	0.19	ng/ul	0.00

Target Compounds				Ovalue		
7) Acenaphthylene	13.76	152	565	0.048	ng/ul#	85
11) Pentachlorophenol	16.47	266	227	0.151	ng/ul	93
12) Phenanthrene	16.85	178	2718	0.156	ng/ul	99
13) Anthracene	16.94	178	1517	0.102	ng/ul	98
15) Fluoranthene	18.88	202	12547	0.607	ng/ul	99
17) Pyrene	19.25	202	13823	0.717	ng/ul	99
18) Benzo(a)anthracene	21.01	228	9841	0.583	ng/ul	99
19) Chrysene	21.06	228	11286	0.605	ng/ul	99
21) Benzo(b)fluoranthene	22.51	252	13654	0.773	ng/ul	94
22) Benzo(k)fluoranthene	22.55	252	8508	0.436	ng/ul	95
23) Benzo(a)pyrene	23.04	252	8976	0.548	ng/ul	92
24) Indeno(1,2,3-cd)pyrene	25.18	276	8438	0.392	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.19	278	4842	0.285	ng/ul	96
26) Benzo(g,h,i)perylene	25.81	276	7073	0.363	ng/ul	97

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

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