

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121018\
 Data File : BN003779.D
 Acq On : 10 Dec 2018 15:22
 Operator : JU/SJ
 Sample : J6155-23DL 2X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4132DL

Quant Time: Dec 11 04:04:13 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 11 03:42:06 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	4154	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	16867	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	8771	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	18815	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	14066	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	12052	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	3536	0.14	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	6997	0.13	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.68	128	15416	0.325	ng/ul	99
7) Acenaphthylene	14.19	152	11605	0.294	ng/ul	99
8) Acenaphthene	14.53	153	9921	0.287	ng/ul	99
9) Fluorene	15.52	166	12851	0.318	ng/ul	100
11) Pentachlorophenol	16.86	266	2342	0.691	ng/ul	99
12) Phenanthrene	17.26	178	17077	0.277	ng/ul	99
13) Anthracene	17.35	178	8623	0.170	ng/ul	99
17) Pyrene	19.64	202	36367	0.721	ng/ul	100
18) Benzo(a)anthracene	21.38	228	12789	0.288	ng/ul	98
19) Chrysene	21.44	228	19809	0.390	ng/ul	100
22) Benzo(k)fluoranthene	23.07	252	8638	0.178	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	26.07	276	12365	0.250	ng/ul#	92
26) Benzo(g,h,i)perylene	26.79	276	14123	0.328	ng/ul	97

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

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