

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN051319\
 Data File : BN005648.D
 Acq On : 13 May 2019 11:55
 Operator : JU/SJ
 Sample : K2606-03
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A5120

Quant Time: May 13 12:28:02 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN050719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun May 12 00:36:09 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	1076	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.36	136	4318	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	2992	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.98	188	7868	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	7969	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	8657	0.40	ng/ul	0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	2305	0.34	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	8742	0.35	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.41	128	7103	0.622	ng/ul	99
5) 2-Methylnaphthalene	12.04	142	4972	0.637	ng/ul	98
9) Fluorene	15.29	166	8818	0.708	ng/ul	99
11) Pentachlorophenol	16.66	266	3800	1.864	ng/ul	92
12) Phenanthrene	17.03	178	17262	0.805	ng/ul	99
15) Fluoranthene	19.06	202	18157	0.612	ng/ul	97
18) Benzo(a)anthracene	21.21	228	23186	0.844	ng/ul	100
22) Benzo(k)fluoranthene	22.83	252	21299	0.653	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.64	276	17001	0.523	ng/ul#	88
26) Benzo(g,h,i)perylene	26.32	276	27658	0.997	ng/ul	96

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

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