

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111918\
 Data File : BN003596.D
 Acq On : 19 Nov 2018 16:20
 Operator : MJ/SJ
 Sample : J5928-11
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A40S5

Quant Time: Nov 19 17:34:52 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 : Mon Nov 19 15:40:59 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	5916	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	21869	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	10408	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	25992	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	27916	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	24907	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	11713	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	30729	0.40	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.69	128	15931	0.259	ng/ul	99
5) 2-Methylnaphthalene	12.30	142	10359	0.242	ng/ul	100
7) Acenaphthylene	14.20	152	38594	0.823	ng/ul	100
8) Acenaphthene	14.54	153	14000	0.341	ng/ul	100
9) Fluorene	15.53	166	17687	0.369	ng/ul	100
11) Pentachlorophenol	16.87	266	535	0.114	ng/ul	97
12) Phenanthrene	17.26	178	63171	0.741	ng/ul	99
13) Anthracene	17.36	178	17126	0.244	ng/ul	99
15) Fluoranthene	19.28	202	42313	0.420	ng/ul	100
17) Pyrene	19.64	202	46302	0.462	ng/ul	100
18) Benzo(a)anthracene	21.38	228	63704	0.722	ng/ul	99
19) Chrysene	21.43	228	41342	0.410	ng/ul	100
21) Benzo(b)fluoranthene	23.01	252	39815	0.386	ng/ul	100
22) Benzo(k)fluoranthene	23.01	252	39866	0.397	ng/ul	100
23) Benzo(a)pyrene	23.61	252	45261	0.507	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	26.06	276	39165	0.383	ng/ul	99
25) Dibenzo(a,h)anthracene	26.08	278	33728	0.410	ng/ul	99
26) Benzo(g,h,i)perylene	26.79	276	40784	0.459	ng/ul	99

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

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