

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN041819\
 Data File : BN005151.D
 Acq On : 18 Apr 2019 13:16
 Operator : JU/SJ
 Sample : K2365-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 X2G53

Quant Time: Apr 19 05:05:26 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN041819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 18 11:57:47 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	2787	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	9952	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	5398	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	12645	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	10807	0.40	ng/ul	0.00
20) Perylene-d12	23.39	264	10901	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	5615	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	13489	0.38	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.42	128	9049	0.329	ng/ul	100
5) 2-Methylnaphthalene	12.04	142	11184	0.573	ng/ul	100
7) Acenaphthylene	13.95	152	6862	0.253	ng/ul	99
9) Fluorene	15.30	166	12714	0.542	ng/ul	98
11) Pentachlorophenol	16.64	266	3347	1.091	ng/ul	99
12) Phenanthrene	17.03	178	13261	0.337	ng/ul	100
13) Anthracene	17.12	178	18769	0.523	ng/ul	99
15) Fluoranthene	19.06	202	13347	0.300	ng/ul	99
17) Pyrene	19.42	202	17096	0.367	ng/ul	99
18) Benzo(a)anthracene	21.19	228	19863	0.482	ng/ul	99
19) Chrysene	21.24	228	21189	0.500	ng/ul	99
21) Benzo(b)fluoranthene	22.74	252	31578	0.752	ng/ul	98
22) Benzo(k)fluoranthene	22.78	252	13768	0.333	ng/ul	99
23) Benzo(a)pyrene	23.29	252	12377	0.332	ng/ul	100
25) Dibenzo(a,h)anthracene	25.58	278	27831	0.823	ng/ul	99
26) Benzo(g,h,i)perylene	26.21	276	12440	0.353	ng/ul	99

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

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