

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN041819\
 Data File : BN005152.D
 Acq On : 18 Apr 2019 15:06
 Operator : JU/SJ
 Sample : K2365-08
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 X2G58

Quant Time: Apr 19 05:15:41 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	3471	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	13324	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	7729	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	18932	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	15723	0.40	ng/ul	0.00
20) Perylene-d12	23.38	264	14597	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	5663	0.26	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	15890	0.30	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.42	128	15148	0.411	ng/ul	99
5) 2-Methylnaphthalene	12.04	142	12258	0.469	ng/ul	100
7) Acenaphthylene	13.95	152	19544	0.504	ng/ul	99
8) Acenaphthene	14.30	153	16592	0.572	ng/ul	100
11) Pentachlorophenol	16.64	266	4595	1.000	ng/ul	99
13) Anthracene	17.12	178	25379	0.472	ng/ul	100
17) Pyrene	19.42	202	45656	0.674	ng/ul	98
19) Chrysene	21.24	228	32106	0.521	ng/ul	99
21) Benzo(b)fluoranthene	22.74	252	33298	0.592	ng/ul	98
22) Benzo(k)fluoranthene	22.78	252	54877	0.993	ng/ul	98
23) Benzo(a)pyrene	23.29	252	22681	0.455	ng/ul	99
25) Dibenzo(a,h)anthracene	25.57	278	27826	0.614	ng/ul	97

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

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