

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN051619\
 Data File : BN005714.D
 Acq On : 16 May 2019 22:02
 Operator : JU/SJ
 Sample : K2690-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :

Quant Time: May 17 04:56:16 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 : Tue May 14 07:15:06 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	1358	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.35	136	5548	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.22	164	3547	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.98	188	8363m	0.40	ng/ul	-0.01
16) Chrysene-d12	21.21	240	5210m	0.40	ng/ul	0.00
20) Perylene-d12	23.42	264	5173m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	2126	0.24	ng/ul	-0.01
14) Fluoranthene-d10	19.02	212	5004m	0.19	ng/ul	-0.01
Target Compounds				Qvalue		
3) Naphthalene	10.39	128	320	0.022	ng/ul#	46
7) Acenaphthylene	13.94	152	545	0.031	ng/ul#	43
9) Fluorene	15.28	166	348	0.024	ng/ul#	69
12) Phenanthrene	17.02	178	12544	0.551	ng/ul	97
15) Fluoranthene	19.05	202	39547m	1.254	ng/ul	
17) Pyrene	19.42	202	34707	1.804	ng/ul	97
18) Benzo(a)anthracene	21.19	228	7465m	0.416	ng/ul	
19) Chrysene	21.24	228	14637m	0.664	ng/ul	
21) Benzo(b)fluoranthene	22.76	252	24886m	1.600	ng/ul	
22) Benzo(k)fluoranthene	22.80	252	6393m	0.328	ng/ul	
23) Benzo(a)pyrene	23.32	252	14781m	0.850	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.62	276	11082m	0.570	ng/ul	
26) Benzo(g,h,i)perylene	26.30	276	18835m	1.136	ng/ul	

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

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