

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121918\
 Data File : BN004121.D
 Acq On : 19 Dec 2018 19:39
 Operator : JU/SJ
 Sample : J6297-11
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E44

Manual Integrations
 APPROVED

Sohil
 12/20/2018 2:03:24 PM

Quant Time: Dec 20 06:18:41 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 : Thu Dec 20 05:49:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	1415	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	7333	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	4491	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	11705m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	12768m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	11246m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	2326	0.21	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	7351m	0.21	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.68	128	1073	0.052	ng/ul#	81
5) 2-Methylnaphthalene	12.29	142	737	0.051	ng/ul	100
12) Phenanthrene	17.26	178	2384m	0.062	ng/ul	
13) Anthracene	17.35	178	1030m	0.033	ng/ul	
15) Fluoranthene	19.27	202	4061m	0.090	ng/ul	
17) Pyrene	19.64	202	6550m	0.143	ng/ul	
18) Benzo(a)anthracene	21.39	228	3869	0.096	ng/ul#	89
19) Chrysene	21.44	228	5085	0.110	ng/ul#	91
21) Benzo(b)fluoranthene	23.02	252	7157m	0.154	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	2514m	0.055	ng/ul	
23) Benzo(a)pyrene	23.62	252	4820	0.120	ng/ul#	76
24) Indeno(1,2,3-cd)pyrene	26.09	276	3989	0.086	ng/ul#	98
26) Benzo(g,h,i)perylene	26.81	276	7660	0.191	ng/ul#	86

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

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