

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN121718\
 Data File : BN004047.D
 Acq On : 17 Dec 2018 21:56
 Operator : JU/SJ
 Sample : J6236-11
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E21

Manual Integrations
 APPROVED

Sohil
 12/18/2018 5:56:23 PM

Quant Time: Dec 18 06:08:37 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 : Tue Dec 18 04:29:07 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	1600	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	8312	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5129	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	13475m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	14740m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	13998m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	2283	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	7864m	0.20	ng/ul	0.00
Target Compounds						Qvalue
3) Naphthalene	10.68	128	589	0.025	ng/ul#	57
5) 2-Methylnaphthalene	12.29	142	357	0.022	ng/ul	95
7) Acenaphthylene	14.19	152	125318	5.426	ng/ul	99
8) Acenaphthene	14.53	153	1581	0.078	ng/ul#	94
9) Fluorene	15.52	166	2469	0.105	ng/ul#	98
12) Phenanthrene	17.26	178	23549m	0.532	ng/ul	
13) Anthracene	17.35	178	106379m	2.927	ng/ul	
15) Fluoranthene	19.28	202	598836m	11.466	ng/ul	
17) Pyrene	19.64	202	646215	12.224	ng/ul	99
18) Benzo(a)anthracene	21.39	228	441716m	9.488	ng/ul	
19) Chrysene	21.44	228	418735m	7.859	ng/ul	
21) Benzo(b)fluoranthene	23.02	252	720598m	12.445	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	245194m	4.346	ng/ul	
23) Benzo(a)pyrene	23.62	252	336056m	6.702	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.09	276	287648m	5.003	ng/ul	
25) Dibenzo(a,h)anthracene	26.09	278	70114m	1.515	ng/ul	
26) Benzo(g,h,i)perylene	26.81	276	217083m	4.344	ng/ul	

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

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