

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN121718\
 Data File : BN004048.D
 Acq On : 17 Dec 2018 22:31
 Operator : JU/SJ
 Sample : J6236-13
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E23

Manual Integrations
 APPROVED

Sohil
 12/18/2018 5:56:26 PM

Quant Time: Dec 18 06:16:45 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	7.83	152	1494	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	7823	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5111	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	13218m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	14717m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	12886m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	3644	0.30	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	12275m	0.32	ng/ul	0.00
Target Compounds				Qvalue		
7) Acenaphthylene	14.19	152	8720	0.379	ng/ul	99
12) Phenanthrene	17.25	178	5817m	0.134	ng/ul	
13) Anthracene	17.35	178	7602m	0.213	ng/ul	
15) Fluoranthene	19.27	202	45251m	0.883	ng/ul	
17) Pyrene	19.64	202	43513m	0.824	ng/ul	
18) Benzo(a)anthracene	21.39	228	25337m	0.545	ng/ul	
19) Chrysene	21.44	228	30402m	0.572	ng/ul	
21) Benzo(b)fluoranthene	23.02	252	51755m	0.971	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	14865m	0.286	ng/ul	
23) Benzo(a)pyrene	23.62	252	20431m	0.443	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.08	276	19348m	0.366	ng/ul	
25) Dibenzo(a,h)anthracene	26.08	278	4865m	0.114	ng/ul	
26) Benzo(g,h,i)perylene	26.81	276	12636m	0.275	ng/ul	

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

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