

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN101018\
 Data File : BN003039.D
 Acq On : 09 Oct 2018 22:26
 Operator : MJ/SJ
 Sample : J5157-09
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AD1

Manual Integrations
 APPROVED

Sohil
 10/10/2018 4:57:50 PM

Quant Time: Oct 10 02:12:33 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN100918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 10 01:34:31 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1767	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	7569	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	4506	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	10123m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.23	240	8453m	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	6817	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	3754	0.32	ng/ul	-0.01
14) Fluoranthene-d10	19.06	212	9496m	0.33	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.44	128	550	0.028	ng/ul#	71
7) Acenaphthylene	13.98	152	761	0.037	ng/ul#	79
12) Phenanthrene	17.06	178	2039m	0.072	ng/ul	
13) Anthracene	17.15	178	1519m	0.057	ng/ul	
15) Fluoranthene	19.09	202	6323m	0.191	ng/ul	
17) Pyrene	19.45	202	10682m	0.325	ng/ul	
18) Benzo(a)anthracene	21.21	228	3765	0.138	ng/ul#	88
19) Chrysene	21.27	228	5653	0.205	ng/ul#	92
21) Benzo(b)fluoranthene	22.78	252	12309m	0.501	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	3708m	0.146	ng/ul	
23) Benzo(a)pyrene	23.34	252	6053	0.275	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.66	276	3124	0.124	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.65	278	940	0.046	ng/ul#	26
26) Benzo(g,h,i)perylene	26.33	276	2636	0.124	ng/ul#	84

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

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