

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101118\
 Data File : BN003096.D
 Acq On : 11 Oct 2018 10:24
 Operator : MJ/SJ
 Sample : J5230-05DL 10X
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A3YZ2DL

Manual Integrations
 APPROVED

Sohil
 10/11/2018 5:45:19 PM

Quant Time: Oct 11 13:44:45 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 : Thu Oct 11 08:10:30 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1368	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.41	136	5569	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	3314	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	7074	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	4758	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	4902	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	215	0.02	ng/ul	0.03
14) Fluoranthene-d10	19.06	212	499	0.02	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.46	128	561	0.039	ng/ul#	58
8) Acenaphthene	14.33	153	592	0.051	ng/ul	96
9) Fluorene	15.33	166	615	0.047	ng/ul#	95
12) Phenanthrene	17.07	178	11679	0.592	ng/ul	100
13) Anthracene	17.16	178	2233	0.120	ng/ul#	89
15) Fluoranthene	19.09	202	14436	0.624	ng/ul	98
17) Pyrene	19.45	202	12506	0.677	ng/ul	99
18) Benzo(a)anthracene	21.22	228	4859m	0.317	ng/ul	
19) Chrysene	21.27	228	4678	0.301	ng/ul	97
21) Benzo(b)fluoranthene	22.79	252	5814m	0.329	ng/ul	
22) Benzo(k)fluoranthene	22.83	252	2107m	0.115	ng/ul	
23) Benzo(a)pyrene	23.35	252	4148	0.262	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.68	276	2972	0.163	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.67	278	787	0.054	ng/ul#	22
26) Benzo(g,h,i)perylene	26.35	276	2783	0.182	ng/ul#	88

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

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