

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110118\
 Data File : BN003389.D
 Acq On : 01 Nov 2018 21:20
 Operator : MJ/SJ
 Sample : J5635-29MSD
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A40H7MSD

Manual Integrations
 APPROVED

Sohil
 11/2/2018 4:14:12 PM

Quant Time: Nov 02 05:19:39 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN110118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 02 04:49:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	11270	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	47397	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	26820	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	65487	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	64545	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	55853	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	21049	0.28	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	55988	0.30	ng/ul	0.00
Target Compounds				Ovalue		
8) Acenaphthene	14.55	153	29392	0.282	ng/ul	98
11) Pentachlorophenol	16.87	266	2860	0.218	ng/ul	99
12) Phenanthrene	17.27	178	49471	0.230	ng/ul	99
13) Anthracene	17.36	178	5843m	0.033	ng/ul	
15) Fluoranthene	19.29	202	135678	0.562	ng/ul	97
17) Pyrene	19.65	202	183947	0.783	ng/ul	100
18) Benzo(a)anthracene	21.39	228	53544	0.265	ng/ul	97
19) Chrysene	21.45	228	67198	0.292	ng/ul	97
21) Benzo(b)fluoranthene	23.02	252	82911m	0.345	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	31677m	0.136	ng/ul	
23) Benzo(a)pyrene	23.62	252	51382	0.256	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	26.08	276	33269	0.155	ng/ul#	99
25) Dibenzo(a,h)anthracene	26.09	278	7885	0.045	ng/ul#	27
26) Benzo(g,h,i)perylene	26.81	276	24989m	0.135	ng/ul	

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

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