

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110118\
 Data File : BN003388.D
 Acq On : 01 Nov 2018 20:45
 Operator : MJ/SJ
 Sample : J5635-28MS
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A40H7MS

Manual Integrations
 APPROVED

Sohil
 11/2/2018 4:14:10 PM

Quant Time: Nov 02 05:17:29 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	11515	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	47812	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	27090	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	66414	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	64207	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	54241	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	21711	0.28	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	57783	0.31	ng/ul	0.00
Target Compounds				Ovalue		
8) Acenaphthene	14.55	153	29130	0.277	ng/ul	99
11) Pentachlorophenol	16.87	266	3078	0.231	ng/ul	99
12) Phenanthrene	17.27	178	38375	0.176	ng/ul	99
13) Anthracene	17.36	178	4418	0.025	ng/ul#	92
15) Fluoranthene	19.29	202	124274	0.508	ng/ul	97
17) Pyrene	19.65	202	172953	0.741	ng/ul	99
18) Benzo(a)anthracene	21.39	228	46614	0.232	ng/ul	97
19) Chrysene	21.45	228	63575	0.278	ng/ul	97
21) Benzo(b)fluoranthene	23.02	252	78704m	0.337	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	30136m	0.133	ng/ul	
23) Benzo(a)pyrene	23.62	252	45348	0.232	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	26.08	276	30233	0.145	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.08	278	7056	0.041	ng/ul#	26
26) Benzo(g,h,i)perylene	26.80	276	22854m	0.128	ng/ul	

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

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