

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE070618\
 Data File : BE096865.D
 Acq On : 6 Jul 2018 22:27
 Operator : SJ/JU
 Sample : J3721-13
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F1H24

Manual Integrations
 APPROVED

Sohil
 7/9/2018 4:14:11 PM

Quant Time: Jul 09 05:32:50 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE070618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 06 16:55:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	3786	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	19840	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	12742	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	30858m	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	30068m	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	27303m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	9138	0.28	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	23889m	0.31	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.19	128	3071	0.064	ng/ul#	95
5) 2-Methylnaphthalene	11.81	142	806	0.025	ng/ul	99
7) Acenaphthylene	13.75	152	8158	0.156	ng/ul	97
8) Acenaphthene	14.09	153	1455	0.034	ng/ul#	93
9) Fluorene	15.08	166	1743	0.039	ng/ul	90
12) Phenanthrene	16.82	178	20228m	0.254	ng/ul	
13) Anthracene	16.91	178	9336m	0.138	ng/ul	
15) Fluoranthene	18.84	202	162683m	1.920	ng/ul	
17) Pyrene	19.21	202	195040	1.976	ng/ul#	80
18) Benzo(a)anthracene	20.96	228	108016m	1.523	ng/ul	
19) Chrysene	21.02	228	159200	1.733	ng/ul	98
21) Benzo(b)fluoranthene	22.54	252	283837m	3.809	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	81830m	1.006	ng/ul	
23) Benzo(a)pyrene	23.12	252	154279m	2.448	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.50	276	88457m	1.352	ng/ul	
25) Dibenzo(a,h)anthracene	25.51	278	20585m	0.396	ng/ul	
26) Benzo(g,h,i)perylene	26.20	276	74509m	1.173	ng/ul	

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

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