

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE070618\
 Data File : BE096860.D
 Acq On : 6 Jul 2018 19:22
 Operator : SJ/JU
 Sample : J3721-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F1H14

Manual Integrations
 APPROVED

Sohil
 7/9/2018 4:14:07 PM

Quant Time: Jul 09 05:28:59 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	3657	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	19293	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	11876	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	30614m	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	27900m	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	27219m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	5679	0.18	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	15150m	0.20	ng/ul	0.00
Target Compounds				Qvalue		
7) Acenaphthylene	13.74	152	2404	0.049	ng/ul	93
12) Phenanthrene	16.82	178	2412m	0.031	ng/ul	
13) Anthracene	16.91	178	2072m	0.031	ng/ul	
15) Fluoranthene	18.85	202	15519m	0.185	ng/ul	
17) Pyrene	19.21	202	27469	0.300	ng/ul#	81
18) Benzo(a)anthracene	20.96	228	18216	0.277	ng/ul#	84
19) Chrysene	21.02	228	18655	0.219	ng/ul	99
21) Benzo(b)fluoranthene	22.54	252	40933m	0.551	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	14270m	0.176	ng/ul	
23) Benzo(a)pyrene	23.11	252	26079m	0.415	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.50	276	12376	0.190	ng/ul#	74
25) Dibenzo(a,h)anthracene	25.52	278	2751	0.053	ng/ul	98
26) Benzo(g,h,i)perylene	26.20	276	11198	0.177	ng/ul#	94

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

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