

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE103118\
 Data File : BE098134.D
 Acq On : 1 Nov 2018 2:20
 Operator : SJ/JU
 Sample : J5701-06
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A40J7

Manual Integrations
 APPROVED

Sohil
 11/1/2018 11:08:14 AM

Quant Time: Nov 01 04:36:47 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE102518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 01 02:11:37 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.87	152	1110	0.40	ng/ul	0.00
2) Naphthalene-d8	10.67	136	5483	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	3789	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	8713m	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	9210m	0.40	ng/ul	0.00
20) Perylene-d12	24.01	264	8985m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	1906	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.31	212	5542m	0.20	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.32	178	1282m	0.064	ng/ul	
15) Fluoranthene	19.35	202	5448m	0.212	ng/ul	
17) Pyrene	19.71	202	4497m	0.209	ng/ul	
18) Benzo(a)anthracene	21.45	228	4195m	0.155	ng/ul	
19) Chrysene	21.50	228	3303m	0.158	ng/ul	
21) Benzo(b)fluoranthene	23.23	252	4903m	0.197	ng/ul	
22) Benzo(k)fluoranthene	23.28	252	1625m	0.067	ng/ul	
23) Benzo(a)pyrene	23.90	252	2899m	0.121	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.70	276	2210m	0.081	ng/ul	
26) Benzo(g,h,i)perylene	27.53	276	2029m	0.088	ng/ul	

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

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