

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE103118\
 Data File : BE098133.D
 Acq On : 1 Nov 2018 1:44
 Operator : SJ/JU
 Sample : J5701-03
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A40E7

Manual Integrations
 APPROVED

Sohil
 11/1/2018 11:08:13 AM

Quant Time: Nov 01 04:31:38 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	1160	0.40	ng/ul	0.00
2) Naphthalene-d8	10.67	136	5797	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	4142	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.28	188	9123m	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	9872m	0.40	ng/ul	0.00
20) Perylene-d12	24.02	264	9735m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	3607	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	10988m	0.37	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.32	178	6877	0.327	ng/ul	98
13) Anthracene	17.42	178	1151m	0.055	ng/ul	
15) Fluoranthene	19.35	202	27535m	1.026	ng/ul	
17) Pyrene	19.71	202	22580m	0.980	ng/ul	
18) Benzo(a)anthracene	21.45	228	20498m	0.706	ng/ul	
19) Chrysene	21.50	228	16348	0.728	ng/ul	97
21) Benzo(b)fluoranthene	23.24	252	22812m	0.846	ng/ul	
22) Benzo(k)fluoranthene	23.28	252	5902m	0.224	ng/ul	
23) Benzo(a)pyrene	23.91	252	13283m	0.512	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.70	276	9512m	0.320	ng/ul	
25) Dibenzo(a,h)anthracene	26.72	278	2486m	0.100	ng/ul	
26) Benzo(g,h,i)perylene	27.53	276	8541m	0.341	ng/ul	

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

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