

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE103118\
 Data File : BE098126.D
 Acq On : 31 Oct 2018 21:21
 Operator : SJ/JU
 Sample : J5657-14
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A40E5

Manual Integrations
 APPROVED

Sohil
 11/1/2018 11:08:07 AM

Quant Time: Nov 01 03:07:44 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	1069	0.40	ng/ul	0.00
2) Naphthalene-d8	10.67	136	5258	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	3693	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.28	188	8214m	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	8732m	0.40	ng/ul	0.00
20) Perylene-d12	24.01	264	8299m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	2712	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.31	212	7789m	0.29	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.33	178	8797	0.464	ng/ul	99
13) Anthracene	17.42	178	1413m	0.075	ng/ul	
15) Fluoranthene	19.34	202	28471m	1.178	ng/ul	
17) Pyrene	19.71	202	25027	1.228	ng/ul	99
18) Benzo(a)anthracene	21.45	228	18351m	0.715	ng/ul	
19) Chrysene	21.50	228	16108	0.811	ng/ul	99
21) Benzo(b)fluoranthene	23.24	252	21430m	0.932	ng/ul	
22) Benzo(k)fluoranthene	23.28	252	7597m	0.338	ng/ul	
23) Benzo(a)pyrene	23.90	252	14209m	0.643	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.70	276	9829m	0.388	ng/ul	
25) Dibenzo(a,h)anthracene	26.72	278	2428m	0.115	ng/ul	
26) Benzo(g,h,i)perylene	27.53	276	7744m	0.363	ng/ul	

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

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