

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE103118\
 Data File : BE098118.D
 Acq On : 31 Oct 2018 16:23
 Operator : SJ/JU
 Sample : J5657-15
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A40E6

Manual Integrations
 APPROVED

Sohil
 11/1/2018 11:08:00 AM

Quant Time: Nov 01 02:07:34 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 01:42:08 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.87	152	928	0.40	ng/ul	0.00
2) Naphthalene-d8	10.67	136	4487	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	3188	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	7403m	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	8111m	0.40	ng/ul	0.00
20) Perylene-d12	24.01	264	7485m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	2061	0.26	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	6127m	0.25	ng/ul	0.00
Target Compounds					Ovalue	
12) Phenanthrene	17.32	178	863m	0.051	ng/ul	
15) Fluoranthene	19.34	202	2327m	0.107	ng/ul	
17) Pyrene	19.71	202	1978	0.104	ng/ul#	86
18) Benzo(a)anthracene	21.45	228	1857m	0.078	ng/ul	
19) Chrysene	21.51	228	1377m	0.075	ng/ul	
21) Benzo(b)fluoranthene	23.24	252	1824m	0.088	ng/ul	
22) Benzo(k)fluoranthene	23.29	252	841m	0.041	ng/ul	
23) Benzo(a)pyrene	23.90	252	1340	0.067	ng/ul#	63
24) Indeno(1,2,3-cd)pyrene	26.71	276	1020	0.045	ng/ul#	96
26) Benzo(g,h,i)perylene	27.53	276	831	0.043	ng/ul#	56

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

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