

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE101018\
 Data File : BE097839.D
 Acq On : 10 Oct 2018 22:05
 Operator : SJ/JU
 Sample : J5183-20
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 JLC33

Manual Integrations
 APPROVED

Sohil
 10/11/2018 5:33:45 PM

Quant Time: Oct 11 02:07:11 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE101018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 11 01:20:33 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	2690	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	13594	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	10457	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	30109m	0.40	ng/ul	0.00
16) Chrysene-d12	21.48	240	35963m	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	32839m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	5913	0.28	ng/ul	0.00
14) Fluoranthene-d10	19.33	212	28664m	0.29	ng/ul	0.00
Target Compounds				Ovalue		
11) Pentachlorophenol	16.95	266	215m	0.078	ng/ul	
12) Phenanthrene	17.34	178	7412	0.103	ng/ul	95
15) Fluoranthene	19.36	202	22447m	0.237	ng/ul	
17) Pyrene	19.72	202	22687m	0.257	ng/ul	
18) Benzo(a)anthracene	21.46	228	8483m	0.085	ng/ul	
19) Chrysene	21.52	228	15333m	0.159	ng/ul	
21) Benzo(b)fluoranthene	23.25	252	24225m	0.285	ng/ul	
22) Benzo(k)fluoranthene	23.30	252	6591m	0.069	ng/ul	
23) Benzo(a)pyrene	23.92	252	11138	0.143	ng/ul#	33
24) Indeno(1,2,3-cd)pyrene	26.72	276	11439	0.130	ng/ul#	100
26) Benzo(g,h,i)perylene	27.54	276	12795	0.154	ng/ul#	63

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

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