

Data Path : Z:\HPCHEM1\BNA E\DATA\BE101017\
 Data File : BE094255.D
 Acq On : 10 Oct 2017 15:32
 Operator : SJ/JU
 Sample : I5522-10
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AD9

Manual Integrations
 APPROVED

Sohil
 10/11/2017 7:19:48 PM

Quant Time: Oct 11 07:21:30 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 11 06:34:16 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1830	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	9561	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	5512	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.24	188	11833	0.40	ng/ul	-0.02
16) Chrysene-d12	21.41	240	11765	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	11630m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	1106	0.07	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	3040	0.08	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.29	178	1753	0.057	ng/ul#	95
15) Fluoranthene	19.30	202	4911	0.121	ng/ul	85
17) Pyrene	19.66	202	5034	0.154	ng/ul#	88
18) Benzo(a)anthracene	21.39	228	4359	0.140	ng/ul#	88
19) Chrysene	21.45	228	5357	0.149	ng/ul#	89
21) Benzo(b)fluoranthene	23.15	252	8883	0.281	ng/ul	93
22) Benzo(k)fluoranthene	23.19	252	3738m	0.102	ng/ul	
23) Benzo(a)pyrene	23.81	252	6901m	0.210	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.57	276	5130m	0.153	ng/ul	
25) Dibenzo(a,h)anthracene	26.58	278	1391m	0.049	ng/ul	
26) Benzo(g,h,i)perylene	27.40	276	4133m	0.143	ng/ul	

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

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