

Data Path : Z:\HPCHEM1\BNA E\DATA\BE101017\
 Data File : BE094254.D
 Acq On : 10 Oct 2017 14:53
 Operator : SJ/JU
 Sample : I5522-09
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AD8

Manual Integrations
 APPROVED

Sohil
 10/11/2017 7:19:46 PM

Quant Time: Oct 11 07:15:44 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	7.91	152	1744	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	9003	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	5163	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.24	188	11038	0.40	ng/ul	-0.02
16) Chrysene-d12	21.41	240	10959	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	11016m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	1529	0.11	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	7982	0.22	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.29	178	2955	0.102	ng/ul	95
15) Fluoranthene	19.30	202	11085	0.292	ng/ul	85
17) Pyrene	19.66	202	11354	0.372	ng/ul#	87
18) Benzo(a)anthracene	21.39	228	5487	0.189	ng/ul#	88
19) Chrysene	21.45	228	7465	0.222	ng/ul	92
21) Benzo(b)fluoranthene	23.15	252	9375	0.313	ng/ul	91
22) Benzo(k)fluoranthene	23.20	252	3614m	0.104	ng/ul	
23) Benzo(a)pyrene	23.81	252	6196m	0.199	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.58	276	4607m	0.145	ng/ul	
25) Dibenzo(a,h)anthracene	26.58	278	1105m	0.041	ng/ul	
26) Benzo(g,h,i)perylene	27.41	276	4480m	0.163	ng/ul	

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

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