

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101417\
 Data File : BE094388.D
 Acq On : 15 Oct 2017 11:59
 Operator : SJ/JU
 Sample : I5548-04DL 5X
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AB5DL

Manual Integrations
 APPROVED

Sohil
 10/16/2017 8:26:02 PM

Quant Time: Oct 16 07:19:29 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 06:50:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	2110	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	11393	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	6898	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	15846	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	15504	0.40	ng/ul	0.00
20) Perylene-d12	23.94	264	14883	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	693	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	1870	0.04	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.29	178	2629	0.063	ng/ul#	91
15) Fluoranthene	19.31	202	7952	0.146	ng/ul	84
17) Pyrene	19.67	202	7091	0.164	ng/ul#	85
18) Benzo(a)anthracene	21.40	228	3552	0.086	ng/ul#	80
19) Chrysene	21.46	228	4860	0.102	ng/ul#	83
21) Benzo(b)fluoranthene	23.17	252	6004m	0.149	ng/ul	
22) Benzo(k)fluoranthene	23.21	252	2112m	0.045	ng/ul	
23) Benzo(a)pyrene	23.82	252	3683	0.088	ng/ul#	77
24) Indeno(1,2,3-cd)pyrene	26.62	276	2779	0.065	ng/ul#	81
26) Benzo(g,h,i)perylene	27.45	276	2326	0.063	ng/ul#	62

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

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