

Data Path : Z:\HPCHEM1\BNA E\DATA\BE112217\
 Data File : BE094835.D
 Acq On : 22 Nov 2017 17:07
 Operator : SJ/JU
 Sample : I6514-09
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AD0

Manual Integrations
 APPROVED

Sohil
 11/27/2017 3:25:11 PM

Quant Time: Nov 23 01:52:48 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE111417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 23 00:56:52 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	752	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.67	136	4893	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.50	164	3493	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.24	188	8523m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.42	240	8767	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	7478m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	2523	0.32	ng/ul	0.01
14) Fluoranthene-d10	19.27	212	7982	0.27	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.27	178	1524m	0.067	ng/ul	
13) Anthracene	17.36	178	774m	0.033	ng/ul	
15) Fluoranthene	19.30	202	10206	0.340	ng/ul	83
17) Pyrene	19.66	202	10308	0.376	ng/ul#	83
18) Benzo(a)anthracene	21.41	228	7046	0.264	ng/ul#	90
19) Chrysene	21.46	228	6548	0.246	ng/ul	96
21) Benzo(b)fluoranthene	23.18	252	6239m	0.280	ng/ul	
22) Benzo(k)fluoranthene	23.22	252	2455m	0.104	ng/ul	
23) Benzo(a)pyrene	23.83	252	3791m	0.171	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.68	276	2107m	0.087	ng/ul	
25) Dibenzo(a,h)anthracene	26.66	278	635m	0.031	ng/ul	
26) Benzo(g,h,i)perylene	27.53	276	1582m	0.078	ng/ul	

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

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