

Data Path : Z:\HPCHEM1\BNA E\DATA\BE032818\
 Data File : BE095886.D
 Acq On : 28 Mar 2018 20:42
 Operator : SJ/JU
 Sample : J1951-05
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Mar 29 05:35:18 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE031318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 29 04:43:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.41	152	1677	0.40	ng/ul	0.00
2) Naphthalene-d8	11.23	136	4581	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.99	164	2963	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.70	188	7362m	0.40	ng/ul	0.00
16) Chrysene-d12	21.79	240	7539m	0.40	ng/ul	0.00
20) Perylene-d12	24.41	264	7916	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.79	152	1532	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.68	212	4761m	0.20	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	11.29	128	1163	0.091	ng/ul#	81
5) 2-Methylnaphthalene	12.87	142	354	0.040	ng/ul	97
8) Acenaphthene	15.06	153	1039	0.092	ng/ul	93
9) Fluorene	16.02	166	765	0.063	ng/ul	91
12) Phenanthrene	17.74	178	12363m	0.561	ng/ul	
13) Anthracene	17.83	178	3580m	0.174	ng/ul	
15) Fluoranthene	19.71	202	28987m	1.085	ng/ul	
17) Pyrene	20.06	202	28996	1.013	ng/ul#	79
18) Benzo(a)anthracene	21.77	228	18671	0.768	ng/ul#	85
19) Chrysene	21.83	228	14717	0.594	ng/ul	99
21) Benzo(b)fluoranthene	23.60	252	21093m	0.769	ng/ul	
22) Benzo(k)fluoranthene	23.64	252	6341m	0.243	ng/ul	
23) Benzo(a)pyrene	24.29	252	15244	0.609	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	27.21	276	9157	0.330	ng/ul#	75
25) Dibenzo(a,h)anthracene	27.22	278	2474	0.111	ng/ul#	96
26) Benzo(a,h,i)perylene	28.09	276	8351	0.350	ng/ul	94

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

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