

Data Path : Z:\HPCHEM1\BNA E\DATA\BE032818\
 Data File : BE095895.D
 Acq On : 29 Mar 2018 2:04
 Operator : SJ/JU
 Sample : J2060-09
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Mar 29 06:41:29 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	1782	0.40	ng/ul	0.00
2) Naphthalene-d8	11.24	136	5119	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.00	164	3461	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.70	188	8069m	0.40	ng/ul	0.00
16) Chrysene-d12	21.79	240	8272m	0.40	ng/ul	0.00
20) Perylene-d12	24.41	264	10095	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.79	152	2991	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.68	212	9577m	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.29	128	5812	0.407	ng/ul#	89
5) 2-Methylnaphthalene	12.87	142	1004	0.101	ng/ul	98
8) Acenaphthene	15.06	153	4739	0.360	ng/ul	97
9) Fluorene	16.02	166	4522	0.317	ng/ul#	91
12) Phenanthrene	17.75	178	77583m	3.211	ng/ul	
13) Anthracene	17.83	178	21579m	0.955	ng/ul	
15) Fluoranthene	19.71	202	126075m	4.307	ng/ul	
17) Pyrene	20.07	202	118093	3.760	ng/ul#	77
18) Benzo(a)anthracene	21.77	228	60300m	2.259	ng/ul	
19) Chrysene	21.83	228	54568	2.008	ng/ul	98
21) Benzo(b)fluoranthene	23.60	252	70604m	2.019	ng/ul	
22) Benzo(k)fluoranthene	23.64	252	21731m	0.654	ng/ul	
23) Benzo(a)pyrene	24.30	252	49171	1.540	ng/ul#	77
24) Indeno(1,2,3-cd)pyrene	27.21	276	28091	0.793	ng/ul#	73
25) Dibenzo(a,h)anthracene	27.22	278	7718	0.273	ng/ul	94
26) Benzo(a,h,i)perylene	28.08	276	25939	0.853	ng/ul#	92

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

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