

Data Path : Z:\HPCHEM1\BNA E\DATA\BE020118\
 Data File : BE095383.D
 Acq On : 2 Feb 2018 6:39
 Operator : SJ/JU
 Sample : J1295-15
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6C13

Manual Integrations
 APPROVED

Sohil
 2/2/2018 4:42:42 PM

Quant Time: Feb 02 07:51:59 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 02 04:53:51 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2069	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	6492	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.01	164	4148	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	8811m	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	8862	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	8188m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	3901	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	10631	0.32	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.74	152	27021	1.354	ng/ul	97
9) Fluorene	16.05	166	2318m	0.145	ng/ul	
11) Pentachlorophenol	17.36	266	31387	18.392	ng/ul	96
12) Phenanthrene	17.75	178	8372	0.283	ng/ul#	90
13) Anthracene	17.84	178	21595	0.796	ng/ul#	93
15) Fluoranthene	19.72	202	308574	8.042	ng/ul	83
17) Pyrene	20.08	202	490369	16.490	ng/ul#	78
18) Benzo(a)anthracene	21.79	228	221388	7.990	ng/ul#	85
19) Chrysene	21.84	228	239411	8.321	ng/ul	96
21) Benzo(b)fluoranthene	23.63	252	324316m	11.227	ng/ul	
22) Benzo(k)fluoranthene	23.67	252	88390m	3.345	ng/ul	
23) Benzo(a)pyrene	24.32	252	76686m	2.973	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.25	276	80762m	2.855	ng/ul	
25) Dibenzo(a,h)anthracene	27.26	278	19881m	0.859	ng/ul	
26) Benzo(g,h,i)perylene	28.12	276	57729m	2.410	ng/ul	

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

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