

Data Path : Z:\HPCHEM1\BNA E\DATA\BE012818\
 Data File : BE095215.D
 Acq On : 28 Jan 2018 20:08
 Operator : SJ/JU
 Sample : J1221-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6A21

Manual Integrations
 APPROVED

Sohil
 1/29/2018 6:25:19 PM

Quant Time: Jan 29 07:33:28 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 29 06:58:46 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	3021	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	9527	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	5969	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	12180	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	13260	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	11690m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	3741	0.24	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	10992	0.24	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.32	128	2742	0.103	ng/ul#	91
5) 2-Methylnaphthalene	12.89	142	1783	0.100	ng/ul	98
7) Acenaphthylene	14.74	152	1134	0.039	ng/ul#	84
8) Acenaphthene	15.08	153	3538	0.159	ng/ul	96
9) Fluorene	16.02	166	2552	0.111	ng/ul#	73
11) Pentachlorophenol	17.36	266	21424	9.081	ng/ul	96
12) Phenanthrene	17.75	178	15419	0.376	ng/ul#	86
13) Anthracene	17.84	178	4606	0.123	ng/ul#	92
15) Fluoranthene	19.72	202	11807	0.223	ng/ul	85
17) Pyrene	20.08	202	9909m	0.223	ng/ul	
18) Benzo(a)anthracene	21.79	228	3442	0.083	ng/ul#	86
19) Chrysene	21.84	228	3581	0.083	ng/ul#	81
21) Benzo(b)fluoranthene	23.62	252	4340m	0.105	ng/ul	
22) Benzo(k)fluoranthene	23.67	252	1446m	0.038	ng/ul	
23) Benzo(a)pyrene	24.32	252	1064m	0.029	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.24	276	1611m	0.040	ng/ul	
26) Benzo(a,h,i)perylene	28.12	276	1578m	0.046	ng/ul	

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

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