

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020318\
 Data File : BE095471.D
 Acq On : 4 Feb 2018 11:15
 Operator : SJ/JU
 Sample : J1315-04
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6C61

Manual Integrations
 APPROVED

Sohil
 2/5/2018 7:01:15 PM

Quant Time: Feb 05 07:35:10 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 Feb 05 04:46:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.43	152	1788	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	5561	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	3693	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	7913	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	7910	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	7367m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	3198	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	8531	0.28	ng/ul	0.00
Target Compounds				Qvalue		
11) Pentachlorophenol	17.36	266	1481	0.966	ng/ul	98
13) Anthracene	17.85	178	1103m	0.045	ng/ul	
15) Fluoranthene	19.72	202	2599	0.075	ng/ul	84
17) Pyrene	20.07	202	2795m	0.105	ng/ul	
18) Benzo(a)anthracene	21.78	228	1012	0.041	ng/ul#	74
19) Chrysene	21.84	228	1057	0.041	ng/ul#	74
21) Benzo(b)fluoranthene	23.62	252	2110m	0.081	ng/ul	
26) Benzo(g,h,i)perylene	28.12	276	573m	0.027	ng/ul	

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

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