

Data Path : Z:\HPCHEM1\BNA E\DATA\BE012718\
 Data File : BE095195.D
 Acq On : 28 Jan 2018 8:03
 Operator : SJ/JU
 Sample : J1220-04
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6A06

Manual Integrations
 APPROVED

Sohil
 1/29/2018 6:22:30 PM

Quant Time: Jan 29 01:14:22 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 00:15:32 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	3527	0.40	ng/ul	0.00
2) Naphthalene-d8	11.27	136	11481	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	7112	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	14704	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	16294	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	17184	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.82	152	3725	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	10926	0.19	ng/ul	0.00
Target Compounds				Ovalue		
11) Pentachlorophenol	17.36	266	1822	0.640	ng/ul	96
12) Phenanthrene	17.75	178	1763	0.036	ng/ul#	92
13) Anthracene	17.85	178	1394	0.031	ng/ul#	91
15) Fluoranthene	19.73	202	2663	0.042	ng/ul	84
17) Pyrene	20.08	202	2522m	0.046	ng/ul	
18) Benzo(a)anthracene	21.79	228	1220	0.024	ng/ul#	67
19) Chrysene	21.84	228	1673	0.032	ng/ul#	72
21) Benzo(b)fluoranthene	23.62	252	2625m	0.043	ng/ul	

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

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