

Data Path : Z:\HPCHEM1\BNA E\DATA\BE012718\
 Data File : BE095184.D
 Acq On : 28 Jan 2018 1:32
 Operator : SJ/JU
 Sample : J1220-03
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6A05

Manual Integrations
 APPROVED

Sohil
 1/29/2018 6:22:04 PM

Quant Time: Jan 28 04:39:18 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jan 27 03:18:43 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	3101	0.40	ng/ul	0.00
2) Naphthalene-d8	11.27	136	9754	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	6183	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.73	188	12975	0.40	ng/ul	0.02
16) Chrysene-d12	21.80	240	14236	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	13450m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.82	152	3492	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	10567	0.21	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.75	152	2975	0.100	ng/ul	97
11) Pentachlorophenol	17.36	266	17941	7.139	ng/ul	95
12) Phenanthrene	17.75	178	3166	0.073	ng/ul#	89
13) Anthracene	17.85	178	5226m	0.131	ng/ul	
15) Fluoranthene	19.73	202	28261	0.500	ng/ul	83
17) Pyrene	20.08	202	29221	0.612	ng/ul#	81
18) Benzo(a)anthracene	21.79	228	10161	0.228	ng/ul#	88
19) Chrysene	21.84	228	14134	0.306	ng/ul	93
21) Benzo(b)fluoranthene	23.62	252	15877m	0.335	ng/ul	
22) Benzo(k)fluoranthene	23.67	252	3564m	0.082	ng/ul	
23) Benzo(a)pyrene	24.32	252	2399m	0.057	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.25	276	3145m	0.068	ng/ul	
26) Benzo(g,h,i)perylene	28.12	276	2666m	0.068	ng/ul	

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

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