

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE071718\
 Data File : BE096979.D
 Acq On : 18 Jul 2018 4:36
 Operator : SJ/JU
 Sample : J3967-04
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 DB1D8

Manual Integrations
 APPROVED

Sohil
 7/18/2018 10:49:59 AM

Quant Time: Jul 18 07:49:38 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE071718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 18 07:15:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	2033	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	10725	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	6928	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	18327m	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	23769m	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	20084m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	5514	0.32	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	18182m	0.33	ng/ul	0.00
Target Compounds				Ovalue		
11) Pentachlorophenol	16.43	266	467m	0.225	ng/ul	
12) Phenanthrene	16.81	178	6562m	0.136	ng/ul	
13) Anthracene	16.90	178	898m	0.021	ng/ul	
15) Fluoranthene	18.84	202	35294m	0.556	ng/ul	
17) Pyrene	19.21	202	30293	0.441	ng/ul#	80
18) Benzo(a)anthracene	20.96	228	7182m	0.114	ng/ul	
19) Chrysene	21.01	228	22911	0.321	ng/ul	97
21) Benzo(b)fluoranthene	22.54	252	37140m	0.652	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	9324m	0.158	ng/ul	
23) Benzo(a)pyrene	23.12	252	15029m	0.300	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.50	276	17751	0.316	ng/ul#	75
25) Dibenzo(a,h)anthracene	25.52	278	3302	0.071	ng/ul#	92
26) Benzo(g,h,i)perylene	26.20	276	19764	0.395	ng/ul	94

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

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