

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080418\
 Data File : BE097217.D
 Acq On : 5 Aug 2018 1:32
 Operator : SJ/JU
 Sample : J4171-21DL 2X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 JJ1L4DL

Manual Integrations
 APPROVED

Sohil
 8/6/2018 12:51:05 PM

Quant Time: Aug 06 04:51:04 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE080118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 06 03:46:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1467	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	7666	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	4648	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	10935m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	11323m	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	10673m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	2007	0.16	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	6029m	0.16	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.73	152	1174	0.058	ng/ul#	89
9) Fluorene	15.08	166	437	0.026	ng/ul#	87
12) Phenanthrene	16.81	178	16967m	0.609	ng/ul	
15) Fluoranthene	18.84	202	23913m	0.649	ng/ul	
17) Pyrene	19.21	202	25012	0.824	ng/ul#	78
18) Benzo(a)anthracene	20.96	228	6368	0.197	ng/ul#	82
19) Chrysene	21.01	228	11312	0.339	ng/ul	97
21) Benzo(b)fluoranthene	22.54	252	11943m	0.411	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	5103m	0.163	ng/ul	
23) Benzo(a)pyrene	23.11	252	9676	0.332	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	25.50	276	7276	0.200	ng/ul#	72
25) Dibenzo(a,h)anthracene	25.52	278	1540	0.052	ng/ul#	68
26) Benzo(g,h,i)perylene	26.20	276	7438	0.236	ng/ul	94

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

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