

Data Path : Z:\HPCHEM1\BNA E\DATA\BE032818\
 Data File : BE095885.D
 Acq On : 28 Mar 2018 20:06
 Operator : SJ/JU
 Sample : J1951-07
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AT0

Manual Integrations
 APPROVED

Sohil
 3/29/2018 5:18:07 PM

Quant Time: Mar 29 05:32:11 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE031318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 29 04:43:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.41	152	1761	0.40	ng/ul	0.00
2) Naphthalene-d8	11.23	136	4773	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.99	164	3039	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.70	188	7541m	0.40	ng/ul	0.00
16) Chrysene-d12	21.79	240	8074m	0.40	ng/ul	0.00
20) Perylene-d12	24.41	264	8686	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.79	152	1454	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.68	212	5121m	0.21	ng/ul	0.00
Target Compounds				Ovalue		
8) Acenaphthene	15.05	153	341	0.030	ng/ul#	88
12) Phenanthrene	17.74	178	6976m	0.309	ng/ul	
13) Anthracene	17.83	178	2075m	0.098	ng/ul	
15) Fluoranthene	19.71	202	29296m	1.071	ng/ul	
17) Pyrene	20.06	202	35756	1.166	ng/ul#	79
18) Benzo(a)anthracene	21.77	228	34183	1.312	ng/ul#	86
19) Chrysene	21.83	228	33513	1.263	ng/ul	99
21) Benzo(b)fluoranthene	23.60	252	44576m	1.482	ng/ul	
22) Benzo(k)fluoranthene	23.64	252	13045m	0.456	ng/ul	
23) Benzo(a)pyrene	24.30	252	22778	0.829	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	27.21	276	11898	0.391	ng/ul#	74
25) Dibenzo(a,h)anthracene	27.22	278	3857	0.158	ng/ul	96
26) Benzo(g,h,i)perylene	28.09	276	6677	0.255	ng/ul	95

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

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