

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022820\
 Data File : BM025154.D
 Acq On : 29 Feb 2020 04:21
 Operator : CG/JU
 Sample : L1506-06
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBCY4

Manual Integrations
 APPROVED

Sohil
 3/3/2020 8:39:26 AM

Quant Time: Mar 02 00:55:07 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM022420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 02 00:36:45 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	2785	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	10688	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	6662	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	15007	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	15328	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	16749	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	4209	0.21	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	11295	0.22	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.03	152	2144	0.056	ng/ul	97
8) Acenaphthene	14.38	153	581	0.023	ng/ul	93
9) Fluorene	15.36	166	770	0.024	ng/ul#	97
12) Phenanthrene	17.10	178	5477	0.104	ng/ul	99
13) Anthracene	17.19	178	6959	0.136	ng/ul	98
15) Fluoranthene	19.11	202	45450	0.593	ng/ul	96
17) Pyrene	19.47	202	57515	0.833	ng/ul	96
18) Benzo(a)anthracene	21.22	228	38760	0.586	ng/ul	100
19) Chrysene	21.27	228	49844	0.756	ng/ul	99
21) Benzo(b)fluoranthene	22.78	252	68456m	1.043	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	26039m	0.400	ng/ul	
23) Benzo(a)pyrene	23.34	252	25983	0.433	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.62	276	15918	0.212	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.62	278	4514	0.072	ng/ul#	77
26) Benzo(g,h,i)perylene	26.28	276	7267	0.115	ng/ul	97

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

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