

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\
 Data File : BM025301.D
 Acq On : 06 Mar 2020 00:22
 Operator : CG/JU
 Sample : L1568-13
 Misc :
 ALS Vial : 98 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBDB4

Manual Integrations
 APPROVED

mohammad
 3/6/2020 10:15:20 PM

Quant Time: Mar 06 03:28:08 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 04 08:02:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	3722	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	16272	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	12121	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	29253m	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	26749	0.40	ng/ul	0.00
20) Perylene-d12	23.42	264	23948	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	7786	0.26	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	25125	0.26	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.49	128	1354	0.028	ng/ul	97
7) Acenaphthylene	14.02	152	10024	0.177	ng/ul	95
8) Acenaphthene	14.36	153	4567	0.104	ng/ul	99
9) Fluorene	15.35	166	5840	0.106	ng/ul	96
12) Phenanthrene	17.08	178	35865	0.380	ng/ul	95
13) Anthracene	17.17	178	41897	0.500	ng/ul	95
15) Fluoranthene	19.10	202	242333	2.053	ng/ul	95
17) Pyrene	19.46	202	250235	2.467	ng/ul	95
18) Benzo(a)anthracene	21.21	228	222972	2.268	ng/ul	99
19) Chrysene	21.26	228	334680	3.098	ng/ul	98
21) Benzo(b)fluoranthene	22.76	252	374519m	3.909	ng/ul	
22) Benzo(k)fluoranthene	22.81	252	121934m	1.213	ng/ul	
23) Benzo(a)pyrene	23.32	252	153653	1.910	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.59	276	86633	0.864	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.60	278	22458	0.270	ng/ul	97
26) Benzo(a,h,i)perylene	26.25	276	55633	0.653	ng/ul#	92

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

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