

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030620\
 Data File : BM025343.D
 Acq On : 07 Mar 2020 11:05
 Operator : CG/JU
 Sample : L1615-19
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBDP3

Manual Integrations
 APPROVED

mohammad
 3/9/2020 3:01:17 PM

Quant Time: Mar 07 21:22:58 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 : Sat Mar 07 19:35:17 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	2134	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	9479	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	6842	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	15758m	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	15191	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	14124	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	4159	0.23	ng/ul	-0.01
14) Fluoranthene-d10	19.07	212	12244	0.24	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.01	152	709	0.022	ng/ul#	69
12) Phenanthrene	17.08	178	2291	0.045	ng/ul	95
13) Anthracene	17.17	178	2633	0.058	ng/ul	97
15) Fluoranthene	19.10	202	3941	0.062	ng/ul	96
17) Pyrene	19.46	202	4477	0.078	ng/ul#	94
18) Benzo(a)anthracene	21.21	228	2453	0.044	ng/ul	91
19) Chrysene	21.26	228	3778	0.062	ng/ul	95
21) Benzo(b)fluoranthene	22.76	252	6983m	0.124	ng/ul	
22) Benzo(k)fluoranthene	22.80	252	2643m	0.045	ng/ul	
23) Benzo(a)pyrene	23.32	252	2356	0.050	ng/ul#	47
24) Indeno(1,2,3-cd)pyrene	25.58	276	2019	0.034	ng/ul#	100
26) Benzo(g,h,i)perylene	26.25	276	1254	0.025	ng/ul#	69

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

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