

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025505.D
 Acq On : 11 Mar 2020 23:33
 Operator : CG/JU
 Sample : L1677-19
 Misc :
 ALS Vial : 98 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0CB6

Manual Integrations
 APPROVED

mohammad
 3/13/2020 8:50:00 AM

Quant Time: Mar 12 02:39:05 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 12 02:03:23 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	3948	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	16612	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	11422	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	23792	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	16487	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	17503	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	5107	0.17	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	11333	0.17	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.47	128	1197	0.024	ng/ul#	85
5) 2-Methylnaphthalene	12.09	142	1009	0.029	ng/ul	97
12) Phenanthrene	17.07	178	11290	0.150	ng/ul	99
13) Anthracene	17.16	178	2213	0.034	ng/ul#	92
15) Fluoranthene	19.09	202	22955	0.281	ng/ul	98
17) Pyrene	19.45	202	20035	0.271	ng/ul	99
18) Benzo(a)anthracene	21.20	228	8494	0.144	ng/ul	96
19) Chrysene	21.25	228	10098	0.153	ng/ul	97
21) Benzo(b)fluoranthene	22.75	252	12917m	0.186	ng/ul	
22) Benzo(k)fluoranthene	22.79	252	4568m	0.063	ng/ul	
23) Benzo(a)pyrene	23.30	252	8251	0.143	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.57	276	6047	0.077	ng/ul#	84
25) Dibenzo(a,h)anthracene	25.58	278	1554	0.024	ng/ul#	48
26) Benzo(g,h,i)perylene	26.23	276	5632	0.083	ng/ul#	91

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

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