

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN080620\
 Data File : BN011375.D
 Acq On : 06 Aug 2020 12:39
 Operator : CG/JU
 Sample : L3419-18DL 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AG7DL

Manual Integrations
 APPROVED

mohammad
 8/7/2020 1:52:58 PM

Quant Time: Aug 06 13:14:17 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 06 12:09:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.46	152	1390	0.40	ng/ul	0.00
2) Naphthalene-d8	10.20	136	4591	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.09	164	2665	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.84	188	5910	0.40	ng/ul	0.00
16) Chrysene-d12	21.06	240	6071	0.40	ng/ul	0.00
20) Perylene-d12	23.18	264	5836	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.82	152	234	0.02	ng/ul	0.01
14) Fluoranthene-d10	18.88	212	604	0.02	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.81	152	1410	0.111	ng/ul#	92
8) Acenaphthene	14.15	153	203	0.020	ng/ul#	93
9) Fluorene	15.15	166	267	0.022	ng/ul#	90
12) Phenanthrene	16.89	178	2765	0.141	ng/ul	96
13) Anthracene	16.98	178	1306	0.076	ng/ul#	87
15) Fluoranthene	18.91	202	9304	0.385	ng/ul	97
17) Pyrene	19.28	202	10003	0.363	ng/ul	99
18) Benzo(a)anthracene	21.05	228	4884	0.209	ng/ul	92
19) Chrysene	21.10	228	12470	0.466	ng/ul	98
21) Benzo(b)fluoranthene	22.56	252	31538m	1.295	ng/ul	
22) Benzo(k)fluoranthene	22.59	252	10644m	0.390	ng/ul	
23) Benzo(a)pyrene	23.09	252	12049	0.569	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.24	276	34239	1.538	ng/ul#	90
25) Dibenzo(a,h)anthracene	25.24	278	7282	0.411	ng/ul	99
26) Benzo(g,h,i)perylene	25.86	276	62447	3.105	ng/ul	97

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

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