

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN011720\
 Data File : BN009395.D
 Acq On : 17 Jan 2020 14:44
 Operator : JU
 Sample : L1109-19
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GASH8

Manual Integrations
 APPROVED

mohammad
 1/21/2020 7:58:04 AM

Quant Time: Jan 17 15:21:56 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN011020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 17 12:28:30 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	1311	0.40	ng/ul	0.00
2) Naphthalene-d8	10.22	136	4523	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.10	164	2551	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.85	188	5510m	0.40	ng/ul	0.00
16) Chrysene-d12	21.06	240	4463	0.40	ng/ul	0.00
20) Perylene-d12	23.18	264	4923	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.82	152	1762	0.24	ng/ul	0.00
14) Fluoranthene-d10	18.89	212	4822	0.29	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.27	128	3100	0.228	ng/ul	96
5) 2-Methylnaphthalene	11.90	142	1909	0.204	ng/ul	99
7) Acenaphthylene	13.82	152	1371	0.105	ng/ul#	90
9) Fluorene	15.16	166	3118	0.259	ng/ul	99
11) Pentachlorophenol	16.54	266	284	0.188	ng/ul	98
12) Phenanthrene	16.89	178	10588	0.556	ng/ul	98
13) Anthracene	16.98	178	864	0.051	ng/ul#	84
15) Fluoranthene	18.92	202	8182	0.363	ng/ul	100
17) Pyrene	19.28	202	9884	0.461	ng/ul	99
18) Benzo(a)anthracene	21.05	228	8140	0.416	ng/ul	100
19) Chrysene	21.10	228	7155	0.338	ng/ul	99
21) Benzo(b)fluoranthene	22.56	252	5045	0.231	ng/ul	92
22) Benzo(k)fluoranthene	22.60	252	9168m	0.368	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.26	276	6904	0.283	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.27	278	5991	0.311	ng/ul#	94
26) Benzo(a,h,i)perylene	25.88	276	9299	0.435	ng/ul	98

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

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