

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101420\
 Data File : BN012217.D
 Acq On : 15 Oct 2020 05:22
 Operator : CG/JU
 Sample : L4184-20
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :

Quant Time: Oct 15 07:31:15 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 15 04:47:44 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	1494	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	6288	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	3759	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	8129m	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	6898	0.40	ng/ul	0.00
20) Perylene-d12	23.39	264	6652	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.99	152	2132	0.23	ng/ul	0.00
14) Fluoranthene-d10	19.04	212	5057	0.23	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	22336	1.231	ng/ul	96
5) 2-Methylnaphthalene	12.07	142	1986	0.172	ng/ul	100
7) Acenaphthylene	13.98	152	3138	0.192	ng/ul#	93
8) Acenaphthene	14.32	153	4229	0.307	ng/ul	98
9) Fluorene	15.31	166	4508	0.294	ng/ul	99
11) Pentachlorophenol	16.66	266	293	0.125	ng/ul	97
12) Phenanthrene	17.05	178	97355	3.757	ng/ul	97
13) Anthracene	17.14	178	16427	0.739	ng/ul	99
15) Fluoranthene	19.07	202	147564	4.870	ng/ul	98
17) Pyrene	19.44	202	123236	3.963	ng/ul	98
18) Benzo(a)anthracene	21.19	228	66130	2.588	ng/ul	98
19) Chrysene	21.24	228	68438	2.199	ng/ul	98
21) Benzo(b)fluoranthene	22.74	252	93438m	3.478	ng/ul	
22) Benzo(k)fluoranthene	22.78	252	26977m	0.907	ng/ul	
23) Benzo(a)pyrene	23.30	252	58854	2.252	ng/ul#	68
24) Indeno(1,2,3-cd)pyrene	25.57	276	46291	1.379	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.58	278	12126	0.451	ng/ul#	82
26) Benzo(g,h,i)perylene	26.24	276	43968	1.429	ng/ul	98

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

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