

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121119\
 Data File : BN008926.D
 Acq On : 11 Dec 2019 20:33
 Operator : JU
 Sample : K6088-11
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BN8

Manual Integrations
 APPROVED

mohammad
 12/12/2019 11:31:50 AM

Quant Time: Dec 12 01:41:00 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 11 10:35:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	2399	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	10535	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	6981	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.93	188	16066	0.40	ng/ul	0.00
16) Chrysene-d12	21.13	240	14412	0.40	ng/ul	0.00
20) Perylene-d12	23.28	264	15466	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	3055	0.18	ng/ul	0.00
14) Fluoranthene-d10	18.96	212	9568	0.20	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.36	128	3869	0.122	ng/ul#	95
5) 2-Methylnaphthalene	11.98	142	7053	0.318	ng/ul	99
7) Acenaphthylene	13.90	152	1086	0.028	ng/ul#	77
9) Fluorene	15.24	166	1303	0.039	ng/ul#	84
11) Pentachlorophenol	16.59	266	147	0.024	ng/ul	97
12) Phenanthrene	16.97	178	29370	0.542	ng/ul	100
13) Anthracene	17.06	178	2113	0.041	ng/ul#	88
15) Fluoranthene	19.00	202	23443	0.352	ng/ul	99
17) Pyrene	19.36	202	21469	0.371	ng/ul	99
18) Benzo(a)anthracene	21.12	228	9772m	0.169	ng/ul	
19) Chrysene	21.17	228	17823	0.313	ng/ul	97
21) Benzo(b)fluoranthene	22.64	252	14071m	0.221	ng/ul	
22) Benzo(k)fluoranthene	22.68	252	4215m	0.067	ng/ul	
23) Benzo(a)pyrene	23.19	252	8063	0.135	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	25.38	276	5517	0.079	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.40	278	1480	0.026	ng/ul#	41
26) Benzo(a,h,i)perylene	26.02	276	5401	0.092	ng/ul	92

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

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