

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121119\
 Data File : BN008959.D
 Acq On : 12 Dec 2019 23:13
 Operator : JU
 Sample : K6089-19
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BR5

Manual Integrations
 APPROVED

mohammad
 12/13/2019 9:08:35 AM

Quant Time: Dec 13 01:44:11 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 : Thu Dec 12 00:49:42 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	2568	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	11058	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.18	164	7085	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.93	188	16245	0.40	ng/ul	0.00
16) Chrysene-d12	21.13	240	13879	0.40	ng/ul	0.00
20) Perylene-d12	23.27	264	13179	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	3317	0.18	ng/ul	0.00
14) Fluoranthene-d10	18.96	212	9667	0.19	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.36	128	8852	0.266	ng/ul	98
5) 2-Methylnaphthalene	11.98	142	6658	0.286	ng/ul	100
7) Acenaphthylene	13.90	152	16928	0.432	ng/ul	99
9) Fluorene	15.24	166	1216	0.036	ng/ul#	79
12) Phenanthrene	16.97	178	26055	0.476	ng/ul	100
13) Anthracene	17.06	178	11170	0.216	ng/ul	98
15) Fluoranthene	18.99	202	88459	1.314	ng/ul	100
17) Pyrene	19.35	202	84045	1.509	ng/ul	99
18) Benzo(a)anthracene	21.11	228	94559	1.694	ng/ul	95
19) Chrysene	21.16	228	113829	2.076	ng/ul	98
21) Benzo(b)fluoranthene	22.64	252	159206m	2.930	ng/ul	
22) Benzo(k)fluoranthene	22.68	252	48377m	0.898	ng/ul	
23) Benzo(a)pyrene	23.18	252	88530	1.736	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	25.38	276	51133	0.857	ng/ul	92
25) Dibenzo(a,h)anthracene	25.38	278	15034	0.312	ng/ul	94
26) Benzo(a,h,i)perylene	26.02	276	43722	0.876	ng/ul	98

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

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