

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121119\
 Data File : BN008938.D
 Acq On : 12 Dec 2019 10:43
 Operator : JU
 Sample : K6088-03DL 5X
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BM9DL

Manual Integrations
 APPROVED

mohammad
 12/13/2019 8:07:38 AM

Quant Time: Dec 12 11:36:37 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	2769	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	11695	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.18	164	7642	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.93	188	17700	0.40	ng/ul	0.00
16) Chrysene-d12	21.13	240	15304	0.40	ng/ul	0.00
20) Perylene-d12	23.28	264	16240	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	757	0.04	ng/ul	0.00
14) Fluoranthene-d10	18.96	212	2295	0.04	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.36	128	8715	0.247	ng/ul	99
5) 2-Methylnaphthalene	11.98	142	6395	0.260	ng/ul	99
7) Acenaphthylene	13.90	152	3282	0.078	ng/ul#	92
8) Acenaphthene	14.24	153	2821	0.090	ng/ul	98
9) Fluorene	15.24	166	4474	0.122	ng/ul#	96
12) Phenanthrene	16.97	178	32456	0.544	ng/ul	100
13) Anthracene	17.06	178	10243	0.182	ng/ul	97
15) Fluoranthene	18.99	202	59920	0.817	ng/ul	99
17) Pyrene	19.35	202	52092	0.848	ng/ul	99
18) Benzo(a)anthracene	21.12	228	32176	0.523	ng/ul	96
19) Chrysene	21.17	228	36903	0.610	ng/ul	96
21) Benzo(b)fluoranthene	22.64	252	36839m	0.550	ng/ul	
22) Benzo(k)fluoranthene	22.68	252	13894m	0.209	ng/ul	
23) Benzo(a)pyrene	23.18	252	21301	0.339	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.38	276	12287	0.167	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.38	278	3459	0.058	ng/ul#	69
26) Benzo(a,h,i)perylene	26.02	276	11177	0.182	ng/ul	98

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

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