

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121319\
 Data File : BN008972.D
 Acq On : 13 Dec 2019 17:47
 Operator : JU
 Sample : K6089-14DL 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BR2DL

Manual Integrations
 APPROVED

mohammad
 12/16/2019 12:23:55 PM

Quant Time: Dec 13 23:02:37 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 : Fri Dec 13 22:56:27 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	13558	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	46603	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.18	164	21067	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.92	188	32424	0.40	ng/ul	0.00
16) Chrysene-d12	21.13	240	14584	0.40	ng/ul	0.00
20) Perylene-d12	23.27	264	16534	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.90	152	6055	0.08	ng/ul	0.00
14) Fluoranthene-d10	18.96	212	7223	0.07	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.36	128	123943	0.882	ng/ul	99
5) 2-Methylnaphthalene	11.98	142	101320	1.033	ng/ul	100
7) Acenaphthylene	13.90	152	11655	0.100	ng/ul#	85
9) Fluorene	15.24	166	3416	0.034	ng/ul#	54
12) Phenanthrene	16.97	178	94196	0.861	ng/ul	99
13) Anthracene	17.06	178	10174	0.098	ng/ul#	95
15) Fluoranthene	18.99	202	65970	0.491	ng/ul	100
17) Pyrene	19.35	202	61215	1.046	ng/ul	99
18) Benzo(a)anthracene	21.11	228	34082	0.581	ng/ul	91
19) Chrysene	21.16	228	45238m	0.785	ng/ul	
21) Benzo(b)fluoranthene	22.64	252	58771m	0.862	ng/ul	
22) Benzo(k)fluoranthene	22.68	252	16999m	0.251	ng/ul	
23) Benzo(a)pyrene	23.18	252	34298	0.536	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.38	276	34640	0.463	ng/ul	90
25) Dibenzo(a,h)anthracene	25.39	278	9688	0.160	ng/ul#	81
26) Benzo(a,h,i)perylene	26.02	276	36951	0.590	ng/ul	98

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

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