

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121319\
 Data File : BN008969.D
 Acq On : 13 Dec 2019 15:50
 Operator : JU
 Sample : K6089-15DL 10X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BR3DL

Manual Integrations
 APPROVED

mohammad
 12/16/2019 12:23:25 PM

Quant Time: Dec 13 17:40:36 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	7.55	152	2595	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	11251	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.18	164	7308	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.92	188	16534	0.40	ng/ul	0.00
16) Chrysene-d12	21.13	240	13830	0.40	ng/ul	0.00
20) Perylene-d12	23.28	264	13068	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	272	0.01	ng/ul	0.00
14) Fluoranthene-d10	18.96	212	913	0.02	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.36	128	9449	0.279	ng/ul	99
5) 2-Methylnaphthalene	11.98	142	8227	0.347	ng/ul	100
7) Acenaphthylene	13.89	152	6480	0.160	ng/ul#	96
9) Fluorene	15.24	166	893	0.025	ng/ul#	77
12) Phenanthrene	16.97	178	29754	0.534	ng/ul	100
13) Anthracene	17.06	178	4529	0.086	ng/ul#	94
15) Fluoranthene	18.99	202	58521	0.854	ng/ul	100
17) Pyrene	19.35	202	53948	0.972	ng/ul	99
18) Benzo(a)anthracene	21.12	228	34995	0.629	ng/ul	95
19) Chrysene	21.17	228	41964	0.768	ng/ul	97
21) Benzo(b)fluoranthene	22.64	252	55077m	1.022	ng/ul	
22) Benzo(k)fluoranthene	22.68	252	21329m	0.399	ng/ul	
23) Benzo(a)pyrene	23.18	252	32220	0.637	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.38	276	19352	0.327	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.39	278	5302	0.111	ng/ul#	83
26) Benzo(a,h,i)perylene	26.02	276	17619	0.356	ng/ul	99

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

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