

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121818\
 Data File : BN004088.D
 Acq On : 18 Dec 2018 21:59
 Operator : JU/SJ
 Sample : J6271-03DL 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9F10DL

Manual Integrations
 APPROVED

Sohil
 12/19/2018 4:16:40 PM

Quant Time: Dec 19 04:21:50 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 19 04:08:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	1682	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	8077	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5013	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	12380	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	13323	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	14846	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	689	0.06	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	2342	0.06	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.68	128	32909	1.451	ng/ul	99
5) 2-Methylnaphthalene	12.29	142	63155	3.987	ng/ul	100
7) Acenaphthylene	14.19	152	1779	0.079	ng/ul#	51
8) Acenaphthene	14.53	153	2569	0.130	ng/ul	90
9) Fluorene	15.52	166	3299	0.143	ng/ul#	91
12) Phenanthrene	17.26	178	329719	8.115	ng/ul	99
13) Anthracene	17.35	178	25833	0.774	ng/ul	98
15) Fluoranthene	19.27	202	59682	1.244	ng/ul	94
17) Pyrene	19.64	202	312987	6.550	ng/ul	97
18) Benzo(a)anthracene	21.39	228	106658m	2.535	ng/ul	
19) Chrysene	21.44	228	193859	4.026	ng/ul	96
21) Benzo(b)fluoranthene	23.02	252	57921m	0.943	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	11089m	0.185	ng/ul	
23) Benzo(a)pyrene	23.62	252	65160	1.225	ng/ul#	95
24) Indeno(1,2,3-cd)pyrene	26.08	276	29667	0.486	ng/ul	97
25) Dibenzo(a,h)anthracene	26.08	278	20110	0.410	ng/ul#	57
26) Benzo(a,h,i)perylene	26.82	276	60827	1.148	ng/ul#	92

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

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