

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111519\
 Data File : BN008589.D
 Acq On : 15 Nov 2019 16:54
 Operator : JU
 Sample : SSTDO.403
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDO.4031

Quant Time: Nov 15 17:27:54 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 Nov 15 17:05:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	2202	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	7255	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.25	164	4241	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	9584	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	10790	0.40	ng/ul	0.00
20) Perylene-d12	23.36	264	10160	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	4777	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.02	212	11890	0.42	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.44	128	8835	0.428	ng/ul	100
5) 2-Methylnaphthalene	12.06	142	6147	0.447	ng/ul	100
7) Acenaphthylene	13.96	152	8826	0.447	ng/ul	100
8) Acenaphthene	14.31	153	6540	0.409	ng/ul	100
9) Fluorene	15.30	166	7735	0.442	ng/ul	100
11) Pentachlorophenol	16.65	266	1449	0.695	ng/ul	100
12) Phenanthrene	17.03	178	13082	0.449	ng/ul	100
13) Anthracene	17.12	178	12236	0.479	ng/ul	100
15) Fluoranthene	19.06	202	15989	0.479	ng/ul	100
17) Pyrene	19.41	202	16397	0.311	ng/ul	100
18) Benzo(a)anthracene	21.17	228	16606	0.442	ng/ul	100
19) Chrysene	21.22	228	16303	0.341	ng/ul	100
21) Benzo(b)fluoranthene	22.72	252	16140	0.426	ng/ul	100
22) Benzo(k)fluoranthene	22.76	252	16531	0.384	ng/ul	100
23) Benzo(a)pyrene	23.27	252	15439	0.405	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.51	276	18004	0.415	ng/ul	100
25) Dibenzo(a,h)anthracene	25.52	278	14530	0.423	ng/ul	100
26) Benzo(g,h,i)perylene	26.15	276	15112	0.389	ng/ul	100

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

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