

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071719\
 Data File : BN006874.D
 Acq On : 17 Jul 2019 16:00
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.432

Quant Time: Jul 17 18:46:07 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN062719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 06:39:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	2227	0.40	ng/ul	0.01
2) Naphthalene-d8	10.37	136	8489	0.40	ng/ul	0.01
6) Acenaphthene-d10	14.25	164	4389	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	9314	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	6907	0.40	ng/ul	0.02
20) Perylene-d12	23.52	264	7555	0.40	ng/ul	0.05
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	4859	0.37	ng/ul	0.01
14) Fluoranthene-d10	19.06	212	9731	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.42	128	9608	0.394	ng/ul	97
5) 2-Methylnaphthalene	12.06	142	5857	0.356	ng/ul	98
7) Acenaphthylene	13.97	152	7956	0.351	ng/ul	99
8) Acenaphthene	14.31	153	6687	0.383	ng/ul	98
9) Fluorene	15.31	166	6951	0.363	ng/ul	98
11) Pentachlorophenol	16.72	266	400	0.288	ng/ul	92
12) Phenanthrene	17.06	178	11147	0.372	ng/ul	99
13) Anthracene	17.16	178	8464	0.327	ng/ul	99
15) Fluoranthene	19.09	202	12698	0.343	ng/ul	99
17) Pyrene	19.46	202	13345	0.423	ng/ul	99
18) Benzo(a)anthracene	21.25	228	9388	0.331	ng/ul	100
19) Chrysene	21.30	228	13572	0.429	ng/ul	99
21) Benzo(b)fluoranthene	22.85	252	11299	0.365	ng/ul	98
22) Benzo(k)fluoranthene	22.89	252	13037	0.402	ng/ul	98
23) Benzo(a)pyrene	23.42	252	11146	0.368	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.76	276	13325	0.434	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.76	278	10099	0.414	ng/ul	97
26) Benzo(g,h,i)perylene	26.44	276	11979	0.467	ng/ul	96

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

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