

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM081920\
 Data File : BM027143.D
 Acq On : 19 Aug 2020 17:12
 Operator : JU/CG
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.429

Quant Time: Aug 19 17:50:13 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM081720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 19 11:15:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	706	0.40	ng/ul	0.00
2) Naphthalene-d8	10.40	136	2238	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	1476	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	3486	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	3427	0.40	ng/ul	0.00
20) Perylene-d12	23.39	264	3078	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	1640	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.04	212	4246	0.43	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.45	128	2801	0.418	ng/ul	98
5) 2-Methylnaphthalene	12.07	142	1962	0.427	ng/ul	98
7) Acenaphthylene	13.98	152	2635	0.411	ng/ul	100
8) Acenaphthene	14.32	153	2350	0.422	ng/ul	99
9) Fluorene	15.31	166	2893	0.433	ng/ul	100
11) Pentachlorophenol	16.67	266	463	0.403	ng/ul	98
12) Phenanthrene	17.05	178	4893	0.415	ng/ul	99
13) Anthracene	17.14	178	4374	0.409	ng/ul	99
15) Fluoranthene	19.07	202	5992	0.423	ng/ul	97
17) Pyrene	19.43	202	5980	0.408	ng/ul#	96
18) Benzo(a)anthracene	21.19	228	5444	0.400	ng/ul	99
19) Chrysene	21.24	228	6487	0.440	ng/ul	98
21) Benzo(b)fluoranthene	22.74	252	6151	0.466	ng/ul	98
22) Benzo(k)fluoranthene	22.78	252	6674	0.486	ng/ul	98
23) Benzo(a)pyrene	23.30	252	5053	0.423	ng/ul#	96
24) Indeno(1,2,3-cd)pyrene	25.58	276	7319	0.444	ng/ul#	91
25) Dibenzo(a,h)anthracene	25.59	278	6059	0.451	ng/ul	97
26) Benzo(g,h,i)perylene	26.24	276	6111	0.429	ng/ul	97

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

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