

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061623\
 Data File : BM040276.D
 Acq On : 16 Jun 2023 12:13
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4121

Quant Time: Jun 16 23:38:10 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM061423.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 15 00:31:16 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.970	152	8349	0.400	ng/ul	0.00
4) Naphthalene-d8	10.778	136	28101	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.609	164	12117	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.353	188	20749	0.400	ng/ul	0.00
17) Chrysene-d12	21.529	240	13215	0.400	ng/ul	# 0.00
23) Perylene-d12	23.952	264	14760	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.350	96	4730	0.459	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.367	152	13312	0.373	ng/ul	0.00
18) Fluoranthene-d10	19.375	212	15510	0.471	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.384	88	4859	0.449	ng/ul	91
5) Naphthalene	10.827	128	26772	0.402	ng/ul	99
7) 2-Methylnaphthalene	12.439	142	15614	0.377	ng/ul	100
8) 1-Methylnaphthalene	12.659	142	15819	0.365	ng/ul	98
10) Acenaphthylene	14.327	152	17752	0.419	ng/ul	100
11) Acenaphthene	14.670	153	15731	0.433	ng/ul	98
12) Fluorene	15.655	166	16184	0.400	ng/ul	98
14) Pentachlorophenol	17.003	266	1473	0.348	ng/ul	99
15) Phenanthrene	17.391	178	22555	0.413	ng/ul	99
16) Anthracene	17.484	178	18491	0.400	ng/ul	99
19) Fluoranthene	19.403	202	20817	0.462	ng/ul	96
20) Pyrene	19.765	202	21414	0.458	ng/ul#	94
21) Benzo(a)anthracene	21.512	228	16332	0.419	ng/ul	99
22) Chrysene	21.564	228	18556	0.450	ng/ul	99
24) Benzo(b)fluoranthene	23.210	252	21802	0.401	ng/ul	94
25) Benzo(k)fluoranthene	23.260	252	20943	0.409	ng/ul#	94
26) Benzo(a)pyrene	23.844	252	18540	0.415	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	26.471	276	27024	0.478	ng/ul#	94
28) Dibenzo(a,h)anthracene	26.487	278	21483	0.474	ng/ul	98
29) Benzo(g,h,i)perylene	27.242	276	23270	0.493	ng/ul	96

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

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