

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070324\
 Data File : BN032439.D
 Acq On : 03 Jul 2024 16:58
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4394

Quant Time: Jul 03 17:44:58 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN062724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 03 03:26:42 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.599	152	4875	0.400	ng/ul	0.00
4) Naphthalene-d8	10.376	136	15942	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.248	164	9764	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.012	188	18666	0.400	ng/ul	0.00
17) Chrysene-d12	21.217	240	13074	0.400	ng/ul	0.00
23) Perylene-d12	23.439	264	15283	0.400	ng/ul	# 0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.143	96	2504	0.462	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.977	152	9599	0.402	ng/ul	0.00
18) Fluoranthene-d10	19.045	212	17635	0.448	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.177	88	2556	0.414	ng/ul#	86
5) Naphthalene	10.426	128	17501	0.415	ng/ul	99
7) 2-Methylnaphthalene	12.054	142	11230	0.402	ng/ul	99
8) 1-Methylnaphthalene	12.274	142	12152	0.419	ng/ul	99
10) Acenaphthylene	13.971	152	18025	0.444	ng/ul	99
11) Acenaphthene	14.313	153	13131	0.417	ng/ul	99
12) Fluorene	15.313	166	14075	0.375	ng/ul	99
14) Pentachlorophenol	16.682	266	1680	0.398	ng/ul	98
15) Phenanthrene	17.054	178	20580	0.402	ng/ul	99
16) Anthracene	17.147	178	19002	0.416	ng/ul	99
19) Fluoranthene	19.078	202	22436	0.451	ng/ul	98
20) Pyrene	19.445	202	22884	0.440	ng/ul	100
21) Benzo(a)anthracene	21.203	228	19290	0.409	ng/ul	98
22) Chrysene	21.255	228	20353	0.393	ng/ul	98
24) Benzo(b)fluoranthene	22.769	252	20609	0.310	ng/ul	89
25) Benzo(k)fluoranthene	22.813	252	21325	0.331	ng/ul#	89
26) Benzo(a)pyrene	23.342	252	20233	0.438	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	25.676	276	26135	0.361	ng/ul	99
28) Dibenzo(a,h)anthracene	25.683	278	20793	0.387	ng/ul#	92
29) Benzo(g,h,i)perylene	26.361	276	21337	0.370	ng/ul	95

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

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