

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070324\
 Data File : BN032474.D
 Acq On : 04 Jul 2024 15:09
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4396

Quant Time: Jul 04 23:06:33 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	5600	0.400	ng/ul	0.00
4) Naphthalene-d8	10.376	136	16836	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.249	164	8867	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.012	188	16250	0.400	ng/ul	0.00
17) Chrysene-d12	21.217	240	14528	0.400	ng/ul	0.00
23) Perylene-d12	23.433	264	17778	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.143	96	2766	0.444	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.977	152	9416	0.373	ng/ul	0.00
18) Fluoranthene-d10	19.045	212	16253	0.372	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.177	88	2995	0.423	ng/ul#	92
5) Naphthalene	10.426	128	18430	0.414	ng/ul	99
7) 2-Methylnaphthalene	12.054	142	10859	0.368	ng/ul	99
8) 1-Methylnaphthalene	12.274	142	11620	0.379	ng/ul	99
10) Acenaphthylene	13.971	152	16382	0.444	ng/ul	99
11) Acenaphthene	14.309	153	11927	0.417	ng/ul	98
12) Fluorene	15.313	166	12428	0.364	ng/ul	98
14) Pentachlorophenol	16.683	266	1534	0.417	ng/ul	98
15) Phenanthrene	17.054	178	17527	0.393	ng/ul	100
16) Anthracene	17.147	178	16344	0.411	ng/ul	98
19) Fluoranthene	19.078	202	20617	0.373	ng/ul	98
20) Pyrene	19.440	202	21652	0.375	ng/ul	98
21) Benzo(a)anthracene	21.203	228	21022	0.401	ng/ul	99
22) Chrysene	21.252	228	22525	0.391	ng/ul	98
24) Benzo(b)fluoranthene	22.766	252	24096	0.312	ng/ul	92
25) Benzo(k)fluoranthene	22.810	252	24178	0.322	ng/ul#	93
26) Benzo(a)pyrene	23.336	252	23283	0.433	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	25.673	276	29919	0.356	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.680	278	23860	0.382	ng/ul	96
29) Benzo(g,h,i)perylene	26.354	276	24461	0.365	ng/ul	96

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

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