

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072624\
 Data File : BN033015.D
 Acq On : 27 Jul 2024 02:03
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4284

Quant Time: Jul 27 03:05:57 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN072524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 25 15:43:00 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.537	152	2949	0.400	ng/ul	0.00
4) Naphthalene-d8	10.306	136	11704	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.166	164	7550	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.945	188	16160	0.400	ng/ul	0.00
17) Chrysene-d12	21.154	240	14404	0.400	ng/ul	0.00
23) Perylene-d12	23.343	264	16022	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.069	96	1611	0.432	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.917	152	6960	0.364	ng/ul	0.00
18) Fluoranthene-d10	18.986	212	14027	0.344	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.098	88	1704	0.401	ng/ul#	84
5) Naphthalene	10.355	128	12439	0.414	ng/ul	96
7) 2-Methylnaphthalene	11.989	142	7743	0.380	ng/ul	98
8) 1-Methylnaphthalene	12.198	142	8834	0.401	ng/ul	99
10) Acenaphthylene	13.893	152	13279	0.430	ng/ul	99
11) Acenaphthene	14.231	153	10329	0.444	ng/ul	99
12) Fluorene	15.249	166	11196	0.413	ng/ul	100
14) Pentachlorophenol	16.641	266	842	0.256	ng/ul	96
15) Phenanthrene	16.987	178	17019	0.393	ng/ul	99
16) Anthracene	17.093	178	14843	0.381	ng/ul	99
19) Fluoranthene	19.014	202	17113	0.361	ng/ul	99
20) Pyrene	19.381	202	18896	0.372	ng/ul	99
21) Benzo(a)anthracene	21.139	228	18732	0.381	ng/ul	98
22) Chrysene	21.189	228	24178	0.422	ng/ul	98
24) Benzo(b)fluoranthene	22.685	252	20303	0.434	ng/ul	92
25) Benzo(k)fluoranthene	22.729	252	24896	0.444	ng/ul#	90
26) Benzo(a)pyrene	23.255	252	20776	0.472	ng/ul#	88
27) Indeno(1,2,3-cd)pyrene	25.554	276	22818	0.341	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.554	278	18184	0.343	ng/ul	95
29) Benzo(g,h,i)perylene	26.224	276	18837	0.354	ng/ul	96

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

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