

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110624\
 Data File : BM048501.D
 Acq On : 07 Nov 2024 07:26
 Operator : RC/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4077

Quant Time: Nov 07 09:54:49 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM110624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 06 14:39:10 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.687	152	3981	0.400	ng/ul	0.00
4) Naphthalene-d8	10.475	136	13820	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.326	164	7879	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.074	188	15140	0.400	ng/ul	0.00
17) Chrysene-d12	21.254	240	12763	0.400	ng/ul	0.00
23) Perylene-d12	23.478	264	14611	0.400	ng/ul	#-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.138	96	1968	0.410	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.075	152	6798	0.380	ng/ul	0.00
18) Fluoranthene-d10	19.100	212	13818	0.386	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.172	88	2308	0.418	ng/ul#	73
5) Naphthalene	10.524	128	13980	0.404	ng/ul	100
7) 2-Methylnaphthalene	12.152	142	8375	0.384	ng/ul	96
8) 1-Methylnaphthalene	12.361	142	9158	0.411	ng/ul	97
10) Acenaphthylene	14.044	152	12418	0.415	ng/ul	98
11) Acenaphthene	14.386	153	9568	0.400	ng/ul	99
12) Fluorene	15.386	166	10231	0.393	ng/ul	99
14) Pentachlorophenol	16.744	266	1542	0.277	ng/ul	97
15) Phenanthrene	17.116	178	15014	0.378	ng/ul	99
16) Anthracene	17.217	178	13716	0.371	ng/ul	99
19) Fluoranthene	19.132	202	18563	0.393	ng/ul	97
20) Pyrene	19.495	202	19872	0.388	ng/ul	98
21) Benzo(a)anthracene	21.239	228	16131	0.386	ng/ul	100
22) Chrysene	21.292	228	20325	0.405	ng/ul	99
24) Benzo(b)fluoranthene	22.808	252	18032	0.396	ng/ul	94
25) Benzo(k)fluoranthene	22.852	252	20831	0.390	ng/ul	94
26) Benzo(a)pyrene	23.381	252	17976	0.404	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	25.728	276	24248	0.391	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.742	278	18765	0.389	ng/ul	94
29) Benzo(g,h,i)perylene	26.413	276	20094	0.386	ng/ul	98

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

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