

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061824\
 Data File : BM046049.D
 Acq On : 18 Jun 2024 13:55
 Operator : MA/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV021

Quant Time: Jun 18 14:27:13 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM061824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 18 13:58:29 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.459	152	2735	0.400	ng/ul	0.00
4) Naphthalene-d8	10.253	136	7529	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.112	164	4057	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.861	188	7866	0.400	ng/ul	0.00
17) Chrysene-d12	21.044	240	7001	0.400	ng/ul	0.00
23) Perylene-d12	23.139	264	10584	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.046	96	1271	0.339	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.881	152	3363	0.333	ng/ul	0.02
18) Fluoranthene-d10	18.889	212	7000	0.334	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.075	88	1390	0.330	ng/ul	96
5) Naphthalene	10.308	128	7455	0.355	ng/ul	99
7) 2-Methylnaphthalene	11.958	142	4104	0.347	ng/ul	100
8) 1-Methylnaphthalene	12.162	142	4269	0.329	ng/ul	98
10) Acenaphthylene	13.839	152	7065	0.367	ng/ul	99
11) Acenaphthene	14.177	153	5107	0.368	ng/ul	99
12) Fluorene	15.176	166	5498	0.369	ng/ul	100
14) Pentachlorophenol	16.566	266	730	0.310	ng/ul	98
15) Phenanthrene	16.904	178	8042	0.364	ng/ul	99
16) Anthracene	17.001	178	6161	0.374	ng/ul	100
19) Fluoranthene	18.917	202	10293	0.340	ng/ul	99
20) Pyrene	19.280	202	11251	0.343	ng/ul	99
21) Benzo(a)anthracene	21.029	228	8707	0.322	ng/ul	99
22) Chrysene	21.079	228	11912	0.345	ng/ul	100
24) Benzo(b)fluoranthene	22.517	252	12954	0.340	ng/ul	98
25) Benzo(k)fluoranthene	22.560	252	14774	0.326	ng/ul	96
26) Benzo(a)pyrene	23.046	252	13133	0.372	ng/ul	91
27) Indeno(1,2,3-cd)pyrene	25.199	276	20180	0.349	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.213	278	15130	0.351	ng/ul	97
29) Benzo(g,h,i)perylene	25.820	276	17256	0.348	ng/ul	98

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

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