

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040424\
 Data File : BM044902.D
 Acq On : 03 Apr 2024 21:51
 Operator : MA/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV028

Quant Time: Apr 03 23:34:53 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM040424.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 03 23:34:03 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.652	152	2436	0.400	ng/ul	0.00
4) Naphthalene-d8	10.435	136	7178	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.302	164	3919	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.064	188	9138	0.400	ng/ul	0.00
17) Chrysene-d12	21.269	240	9214	0.400	ng/ul	0.00
23) Perylene-d12	23.502	264	11847	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.226	96	1188	0.371	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.046	152	3926	0.404	ng/ul	0.00
18) Fluoranthene-d10	19.103	212	9400	0.404	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.256	88	1049	0.342	ng/ul	99
5) Naphthalene	10.485	128	8081	0.424	ng/ul	97
7) 2-Methylnaphthalene	12.123	142	5049	0.444	ng/ul	100
8) 1-Methylnaphthalene	12.332	142	4793	0.386	ng/ul	99
10) Acenaphthylene	14.025	152	8163	0.417	ng/ul	98
11) Acenaphthene	14.367	153	5754	0.411	ng/ul	99
12) Fluorene	15.366	166	6666	0.425	ng/ul	100
14) Pentachlorophenol	16.731	266	792	0.370	ng/ul	96
15) Phenanthrene	17.107	178	10691	0.416	ng/ul	99
16) Anthracene	17.208	178	10283	0.406	ng/ul	99
19) Fluoranthene	19.131	202	13520	0.411	ng/ul	98
20) Pyrene	19.498	202	14392	0.414	ng/ul	100
21) Benzo(a)anthracene	21.257	228	13645	0.425	ng/ul	99
22) Chrysene	21.307	228	16162	0.404	ng/ul	99
24) Benzo(b)fluoranthene	22.830	252	17409	0.407	ng/ul	95
25) Benzo(k)fluoranthene	22.874	252	19743	0.412	ng/ul	97
26) Benzo(a)pyrene	23.406	252	16832	0.412	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.749	276	23541	0.420	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.776	278	18352	0.414	ng/ul	98
29) Benzo(g,h,i)perylene	26.437	276	19212	0.398	ng/ul	99

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

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