

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082423\
 Data File : BM041464.D
 Acq On : 24 Aug 2023 14:15
 Operator : MA/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV049

Quant Time: Aug 24 18:36:23 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM082423.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 24 18:34:12 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.118	152	9388	0.400	ng/ul	0.00
4) Naphthalene-d8	10.950	136	25651	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.759	164	12115	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.515	188	23767	0.400	ng/ul	0.00
17) Chrysene-d12	21.708	240	7604	0.400	ng/ul	0.00
23) Perylene-d12	24.284	264	6551	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.448	96	4034	0.340	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.528	152	11740	0.347	ng/ul	0.00
18) Fluoranthene-d10	19.538	212	13446	0.401	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.486	88	4326	0.354	ng/ul	94
5) Naphthalene	10.999	128	26286	0.344	ng/ul	100
7) 2-Methylnaphthalene	12.599	142	15778	0.344	ng/ul	99
8) 1-Methylnaphthalene	12.814	142	16012	0.343	ng/ul	99
10) Acenaphthylene	14.486	152	24202	0.333	ng/ul	100
11) Acenaphthene	14.823	153	17391	0.344	ng/ul	98
12) Fluorene	15.809	166	19595	0.346	ng/ul	99
14) Pentachlorophenol	17.135	266	1326	0.296	ng/ul	98
15) Phenanthrene	17.557	178	30446	0.349	ng/ul	100
16) Anthracene	17.650	178	26294	0.331	ng/ul	100
19) Fluoranthene	19.566	202	29183	0.418	ng/ul	99
20) Pyrene	19.933	202	29160	0.414	ng/ul	99
21) Benzo(a)anthracene	21.691	228	17911	0.362	ng/ul	99
22) Chrysene	21.746	228	20383	0.381	ng/ul	100
24) Benzo(b)fluoranthene	23.486	252	17688	0.394	ng/ul	98
25) Benzo(k)fluoranthene	23.538	252	17747	0.395	ng/ul	98
26) Benzo(a)pyrene	24.167	252	14077	0.382	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	27.011	276	16141	0.350	ng/ul#	97
28) Dibenzo(a,h)anthracene	27.048	278	12230	0.347	ng/ul	99
29) Benzo(g,h,i)perylene	27.853	276	14523	0.358	ng/ul	98

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

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