

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091523\
 Data File : BM041887.D
 Acq On : 15 Sep 2023 15:37
 Operator : MA/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV035

Quant Time: Sep 16 03:50:21 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM091523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 16 03:49:33 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.971	152	8866	0.400	ng/ul	0.00
4) Naphthalene-d8	10.790	136	22434	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.615	164	11966	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.371	188	25652	0.400	ng/ul	0.00
17) Chrysene-d12	21.542	240	15706	0.400	ng/ul	0.00
23) Perylene-d12	23.962	264	16843	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	4274	0.411	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.368	152	13261	0.431	ng/ul	0.00
18) Fluoranthene-d10	19.389	212	24145	0.417	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.418	88	4586	0.424	ng/ul	93
5) Naphthalene	10.839	128	31480	0.426	ng/ul	100
7) 2-Methylnaphthalene	12.445	142	20097	0.429	ng/ul	99
8) 1-Methylnaphthalene	12.665	142	20457	0.429	ng/ul	100
10) Acenaphthylene	14.342	152	33925	0.413	ng/ul	99
11) Acenaphthene	14.680	153	23794	0.419	ng/ul	100
12) Fluorene	15.665	166	27724	0.418	ng/ul	98
14) Pentachlorophenol	16.995	266	5885	0.386	ng/ul	100
15) Phenanthrene	17.409	178	43877	0.422	ng/ul	99
16) Anthracene	17.502	178	41469	0.418	ng/ul	99
19) Fluoranthene	19.422	202	49553	0.420	ng/ul	100
20) Pyrene	19.789	202	51744	0.419	ng/ul	99
21) Benzo(a)anthracene	21.524	228	40529	0.415	ng/ul	98
22) Chrysene	21.580	228	38270	0.418	ng/ul	99
24) Benzo(b)fluoranthene	23.213	252	39942	0.417	ng/ul	95
25) Benzo(k)fluoranthene	23.263	252	38462	0.418	ng/ul	94
26) Benzo(a)pyrene	23.851	252	33630	0.417	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	26.468	276	43704	0.409	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.485	278	32624	0.410	ng/ul	97
29) Benzo(g,h,i)perylene	27.253	276	35545	0.407	ng/ul	98

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

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