

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090321\
 Data File : BM032043.D
 Acq On : 03 Sep 2021 16:41
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV028

Quant Time: Sep 03 17:15:43 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM090321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 03 16:21:27 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.459	152	1407	0.400	ng/ul	0.00
4) Naphthalene-d8	10.216	136	4927	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.102	164	3141	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.865	188	5769	0.400	ng/ul	0.00
17) Chrysene-d12	21.073	240	3807	0.400	ng/ul	0.00
23) Perylene-d12	23.195	264	3412	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.068	96	609	0.394	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.817	152	3158	0.414	ng/ul	0.00
18) Fluoranthene-d10	18.903	212	4947	0.443	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.102	88	621	0.411	ng/ul	96
5) Naphthalene	10.265	128	5289	0.405	ng/ul	96
7) 2-Methylnaphthalene	11.898	142	3738	0.410	ng/ul	99
8) 1-Methylnaphthalene	12.114	142	3665	0.408	ng/ul	99
10) Acenaphthylene	13.827	152	5095	0.402	ng/ul	99
11) Acenaphthene	14.166	153	4079	0.393	ng/ul	99
12) Fluorene	15.167	166	4714	0.390	ng/ul	100
14) Pentachlorophenol	16.539	266	555	0.364	ng/ul	93
15) Phenanthrene	16.907	178	7022	0.400	ng/ul	99
16) Anthracene	17.007	178	6179	0.401	ng/ul	99
19) Fluoranthene	18.934	202	7202	0.438	ng/ul	99
20) Pyrene	19.300	202	7157	0.422	ng/ul	99
21) Benzo(a)anthracene	21.061	228	5347	0.384	ng/ul	97
22) Chrysene	21.110	228	6340	0.397	ng/ul	99
24) Benzo(b)fluoranthene	22.565	252	6291	0.401	ng/ul	99
25) Benzo(k)fluoranthene	22.608	252	6690	0.413	ng/ul	97
26) Benzo(a)pyrene	23.107	252	5454	0.398	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	25.273	276	7464	0.400	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.290	278	5993	0.399	ng/ul	97
29) Benzo(g,h,i)perylene	25.908	276	6676	0.402	ng/ul	99

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

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