

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121522\
 Data File : BM038029.D
 Acq On : 15 Dec 2022 13:22
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV021

Quant Time: Dec 15 21:57:01 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM121522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 15 21:53:17 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.064	152	3936	0.400	ng/u1	0.00
4) Naphthalene-d8	10.878	136	11561	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.685	164	6174	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.422	188	13043	0.400	ng/u1	0.00
17) Chrysene-d12	21.586	240	8240	0.400	ng/u1	0.00
23) Perylene-d12	24.038	264	7808	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.410	96	1920	0.357	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.456	152	6191	0.367	ng/u1	0.00
18) Fluoranthene-d10	19.436	212	11481	0.384	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.448	88	1874	0.386	ng/u1	98
5) Naphthalene	10.928	128	14154	0.404	ng/u1	100
7) 2-Methylnaphthalene	12.528	142	9057	0.421	ng/u1	99
8) 1-Methylnaphthalene	12.748	142	8776	0.398	ng/u1	100
10) Acenaphthylene	14.407	152	13331	0.412	ng/u1	99
11) Acenaphthene	14.745	153	10014	0.408	ng/u1	99
12) Fluorene	15.730	166	11165	0.411	ng/u1	99
14) Pentachlorophenol	17.075	266	1748	0.324	ng/u1	99
15) Phenanthrene	17.464	178	17277	0.393	ng/u1	99
16) Anthracene	17.552	178	15821	0.390	ng/u1	99
19) Fluoranthene	19.469	202	17863	0.420	ng/u1	99
20) Pyrene	19.831	202	18117	0.417	ng/u1	99
21) Benzo(a)anthracene	21.568	228	14600	0.410	ng/u1	99
22) Chrysene	21.621	228	14918	0.415	ng/u1	100
24) Benzo(b)fluoranthene	23.287	252	14721	0.398	ng/u1	99
25) Benzo(k)fluoranthene	23.336	252	14149	0.396	ng/u1	99
26) Benzo(a)pyrene	23.930	252	12090	0.388	ng/u1	99
27) Indeno(1,2,3-cd)pyrene	26.609	276	13946	0.356	ng/u1#	100
28) Dibenzo(a,h)anthracene	26.622	278	11513	0.378	ng/u1	100
29) Benzo(g,h,i)perylene	27.397	276	12517	0.376	ng/u1	99

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

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